

Maximum k -Plex Search: An Alternated Reduction-and-Bound Method

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ABSTRACT

k -plexes relax cliques by allowing each vertex to disconnect to at most k vertices. Finding a maximum k -plex in a graph is a fundamental operator in graph mining and has been receiving significant attention from various domains. The state-of-the-art algorithms all adopt the branch-reduction-and-bound (BRB) framework where a key step, called *reduction-and-bound* (RB), is used for narrowing down the search space. A common practice of RB in existing works is SeqRB, which *sequentially* conducts the reduction process followed by the bounding process *once* at a branch. However, these algorithms suffer from the efficiency issues. In this paper, we propose a new *alternated reduction-and-bound* method AltRB for conducting RB. AltRB first partitions a branch into two parts and then *alternatively* and *iteratively* conducts the reduction process and the bounding process at each part of a branch. With newly-designed reduction rules and bounding methods, AltRB is superior to SeqRB in effectively narrowing down the search space *in both theory and practice*. Further, to boost the performance of BRB algorithms, we develop efficient and effective pre-processing methods which reduce the size of the input graph and heuristically compute a large k -plex as the lower bound. We conduct extensive experiments on 664 real and synthetic graphs. The experimental results show that our proposed algorithm kPEX with AltRB and novel pre-processing techniques runs up to two orders of magnitude faster and solves more instances than state-of-the-art algorithms.

PVLDB Reference Format:

Shuohao Gao, Kaiqiang Yu, Shengxin Liu, and Cheng Long. Maximum k -Plex Search: An Alternated Reduction-and-Bound Method. PVLDB, 14(1): XXX-XXX, 2020.
doi:XX.XX/XXX.XX

PVLDB Artifact Availability:

The source code, data, and/or other artifacts have been made available at <https://github.com/ShuohaoGao/kPEX>.

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Proceedings of the VLDB Endowment, Vol. 14, No. 1 ISSN 2150-8097.
doi:XX.XX/XXX.XX

1 INTRODUCTION

The graph model serves as a versatile tool for abstracting numerous real-world data which captures relationships between diverse entities in social networks, biological networks, publication networks, and so on. Cohesive subgraph mining is one of the central topics in graph analysis and data mining where the objective is to mine those *dense or cohesive* subgraphs that normally bring valuable insights for analysis [10, 25, 26, 34, 39]. For example, cohesive subgraph mining has been used to detect a terrorist cell in social networks [38], to identify protein complexes in biological networks [62], and to find a group of research collaborators in publication networks [32].

The *clique* is arguably the most well-known cohesive subgraph where every pair of distinct vertices is connected by an edge. In the literature, the study of efficient algorithms for extracting the maximum clique or enumerating maximal cliques is extensive, e.g., [7, 8, 18, 24, 44, 45, 49, 50]. Nevertheless, clique, being a tightly interconnected subgraph, is over-restrictive, which limits its practical usefulness. To circumvent this issue, relaxations of clique have been proposed and studied in the literature, such as k -plex [48], k -core [47], quasi-clique [31, 58], and k -defective clique [9, 21]. In particular, k -plex relaxes clique by allowing each vertex to disconnect to at most k vertices (including the vertex itself). It is clear that 1-plex corresponds to clique. The research of cohesive subgraph mining in the context of k -plex has recently attracted increasing interests [11, 12, 22, 27, 35, 36, 53, 54, 56, 63].

In this paper, we study the *maximum k -plex search* problem which aims to search the k -plex with the largest number of vertices in the given graph. It is well-known that the maximum k -plex search problem is NP-hard for any fixed k [3]. Thus, existing studies and ours focus on designing practically efficient algorithms.

Existing algorithms. The state-of-the-art algorithms all (conceptually) adopt the *branch-reduction-and-bound* (BRB) framework [11, 12, 27, 35, 36, 53, 63]. The idea is to recursively solve the problem instance (or branch) by solving the subproblem instances (or sub-branches) produced via a process of *branching*. A branch denoted by (S, C) corresponds to a problem instance of finding the largest k -plex from the subgraph (of the input graph) induced by vertex set $S \cup C$, where the partial solution S corresponds to a k -plex and the candidate set C corresponds to the set of vertices used to expand the partial solution. We refer the search space of a branch to the set of possible k -plexes in the subgraph induced by $S \cup C$. At each

branch, a key step, named *reduction-and-bound* (RB), is performed for narrowing down the search space. We note that existing studies all follow a sequential framework, called SeqRB, for implementing the RB step. Specifically, SeqRB *sequentially* conducts two processes *once*: 1) the reduction process shrinks the candidate set C by removing some unpromising vertices that cannot appear in the largest k -plex; and 2) the bounding process computes the upper bound of the size of the largest k -plex in the branch refined by the first step, which is used for pruning unnecessary branches (i.e., with the upper bound no larger than the largest k -plex seen so far). The rationale behind is that with some vertices being removed by the reduction process, the bounding process may obtain a smaller upper bound so as to prune more branches.

Existing studies focus solely on sharpening the reduction rules and upper bound computation methods used in SeqRB while devoting little effort to improving the whole RB framework. We observe that, in SeqRB, the reduction process benefits the bounding process, but not the other way; and thus they are sequentially conducted only once. One interesting question is that: *Can we design a new RB framework where the reduction process and the bounding process can benefit each other?*

We remark that some recent studies [11, 12, 35, 53, 63] boost the practical performance of BRB algorithms by devising various pre-processing techniques. These techniques include 1) graph reduction algorithms [11, 63] for reducing the size of the input graph (among which the best one is CTCP [11]); and 2) heuristic algorithms [11, 12, 63] for computing an initial large k -plex used for the above-mentioned reduction algorithms (among which the best ones are kPlex-Degen [11] and EGo-Degen [12]).

Our new methods. In this paper, we first propose a new framework, called *alternated reduction-and-bound* (AltRB), for conducting the RB step at a branch (S, C) . AltRB differs from SeqRB mainly in the way of conducting the reduction process and the bounding process. Specifically, AltRB first partitions a branch into two parts (i.e., $S = S_L \cup S_R$ and $C = C_L \cup C_R$). With newly-designed reduction rules and upper bound computation methods on each part, *the bounding process on one part will benefit the reduction process on the other* (note that the reduction process still benefits the bounding process on the same part, which is the same as SeqRB). Thus, AltRB *alternatively* and *iteratively* conducts the reduction process and the bounding process at each part of a branch (e.g., bounding on $S_L \cup C_L \rightarrow$ reduction on $S_R \cup C_R \rightarrow$ bounding on $S_R \cup C_R \rightarrow$ reduction on $S_L \cup C_L \rightarrow \dots$). In this manner, the bounding process and the reduction process could mutually benefit from each other. We show that AltRB is superior to SeqRB in narrowing down the search space in both theory (as will be shown in Equation (9)) and practice (as will be shown in Table 4). We further design efficient pre-processing techniques for boosting the practical performance of BRB algorithms: 1) a new method CF-CTCP, which differs with CTCP in the way of conducting different reductions at each iteration, and 2) a heuristic algorithm KPHeur is that iteratively compute a large initial maximal k -plex.

With all the above newly-designed techniques, we develop a new BRB algorithm called kPEX, which runs up to two orders of magnitude faster and solves more instances than state-of-the-art algorithms kPlexT [12], kPlexS [11], KPLEX [53], and Di seMKP [35].

Our contributions. Our main contributions are as follows.

- We propose a new BRB algorithm called kPEX, which incorporates the proposed *alternated reduction-and-bound* method AltRB (Section 3). With our newly devised reduction rules and bounding methods, AltRB is superior to SeqRB in narrowing down the search space (Section 4).
- We design efficient pre-processing techniques for boosting the performance of BRB algorithms, namely a new method CF-CTCP for reducing the size of the input graph and a heuristic KPHeur is for computing a large initial k -plex (Section 5).
- We conduct extensive experiments on 664 real and synthetic graphs to verify the effectiveness and efficiency of our algorithms. Compared with the state-of-the-art algorithms, our kPEX 1) solves most number of graph instances within the time limit and 2) runs up to two orders of magnitude faster than existing algorithms (Section 6).

2 PRELIMINARIES

Let $G = (V, E)$ be a simple graph with $|V| = n$ vertices and $|E| = m$ edges. A vertex v is said to be a neighbor of (or adjacent to) vertex u if there is an edge between u and v , i.e., $(u, v) \in E$. Denote by $N_G(u) = \{v \in V \mid (u, v) \in E\}$ and $d_G(u) = |N_G(u)|$ the neighbor set and the degree of the vertex u in G , respectively. Given a vertex subset $S \subseteq V$, we use $G[S]$ to denote the subgraph induced by S , i.e., $G[S] = (S, \{(u, v) \in E \mid u, v \in S\})$, and use $N_G(u, S)$ (resp. $\bar{N}_G(u, S)$) to denote sets of neighbors (resp. non-neighbors that include u itself) of u in $G[S]$. We omit the subscript G when the context is clear. Given a graph g , we use $V(g)$ and $E(g)$ to denote the sets of vertices and edges in g , respectively.

In this paper, we focus on the cohesive subgraph of k -plex.

Definition 2.1 (k -plex [48]). Given a positive integer k , a graph g is said to be a k -plex if $d_g(u) \geq |V(g)| - k$ for each vertex $u \in V(g)$.

Obviously, a 1-plex is a clique where each two vertices are adjacent. Note also that k -plex has the *hereditary* property, i.e., any induced subgraph of a k -plex is also a k -plex [48].

Problem statement. Given a graph $G = (V, E)$ and an integer $k \geq 2$, the *maximum k -plex search problem* aims to find the largest k -plex $G[S]$ with $|S| \geq 2k - 1$ in G .

Following the previous studies [11, 53], we focus on finding k -plexes with at least $2k - 1$ vertices for the following considerations. **First**, the value of k is usually small in real applications, e.g., $k \leq 6$ in [27, 36, 56, 63]. Hence, a k -plex with at most $2k - 2$ vertices is less informative in practice. **Second**, a k -plex with at least $2k - 1$ vertices has the diameter of at most 2 [63], which is more cohesive.

We next introduce some useful concepts used in this paper.

k -core/ k -truss. We review useful cohesive subgraph definitions.

Definition 2.2. Given a positive integer k , a graph g is said to be

- a k -core if $d_g(u) \geq k$ for each vertex $u \in V(g)$ [47];
- a k -truss if each edge $(u, v) \in E(g)$ belongs to at least $k - 2$ triangles, i.e., $|N_g(u) \cap N_g(v)| \geq k - 2$ for each edge $(u, v) \in E(g)$ [16].

Clearly, a k -core g is a $(|V(g)| - k)$ -plex and a k -truss g' is a $(|V(g')| - k + 1)$ -plex.

Algorithm 1: Our framework: kPEX

Input: A graph $G = (V, E)$ and an integer k
Output: The largest k -plex $G[S^*]$
/* Stage-I.1: Heuristic&Preprocessing (Sec. 5) */
1 $S^* \leftarrow$ a large k -plex via a heuristic process KPHeuris;
2 $G \leftarrow$ apply reduction method CF-CTCP to reduce G ;
/* Stage-I.2: Divide-and-conquer framework */
3 **while** $V(G) \neq \emptyset$ **do**
4 $v \leftarrow$ the vertex with the minimum degree in G ;
5 $g \leftarrow$ the subgraph of G induced by $N^{\leq 2}(v)$;
/* Stage-II: branch-reduction-bound (Sec. 4) */
6 BRB_Rec($g, \{v\}, V(g) \setminus \{v\}, k$);
7 $G \leftarrow$ apply reduction method CF-CTCP to reduce G ;
8 **return** $G[S^*]$;
9 **Procedure** BRB_Rec(G, S, C, k)
10 $C^*, UB^* \leftarrow$ AltRB(G, S, C, k);
11 **if** $UB^* \leq |S^*|$ **then return**;
12 **if** $S \cup C^*$ is a k -plex **then** update S^* by $S \cup C^*$ and **return**;
13 $v^* \leftarrow$ a branching vertex selected from C^* ;
14 BRB_Rec($G, S \cup \{v^*\}, C^* \setminus \{v^*\}, k$);
15 BRB_Rec($G, S, C^* \setminus \{v^*\}, k$);

Degeneracy order. The sequence of vertices v_1, v_2, \dots, v_n in $G = (V, E)$ is called the *degeneracy order* of G if v_i has the minimum degree in $G[\{v_i, v_{i+1}, \dots, v_n\}]$ for each v_i in V [4]. Further, the *degeneracy* of G , denoted by $\delta(G)$ (or δ if the context is clear), is defined as the smallest number such that every induced subgraph of G has a vertex of degree at most $\delta(G)$. In other words, G does not have an induced subgraph that is a $(\delta + 1)$ -core. The degeneracy order and the value of δ can be obtained by iteratively peeling the vertex with minimum degree in the current induced subgraph with time complexity of $O(m)$ [4]. Also, it is known that $\delta \leq \sqrt{n + 2m}$ [8].

3 THE FRAMEWORK OF KPEX

Our algorithm, named kPEX, follows the *branch-reduction-and-bound* (BRB) framework which is (conceptually) adopted by existing algorithms [11, 12, 27, 35, 36, 53, 63]. The idea is to recursively partition the current problem instance of finding the largest k -plex into two subproblem instances via a process of *branching*. Specifically, a problem instance (or branch) is denoted by (G, S, C) (or, simply (S, C) when the context is clear) where the *partial solution* S induces a k -plex (i.e., $G[S]$) and the *candidate set* C is a set of vertices that will be used to expand S . Solving the branch (S, C) refers to finding the largest k -plex $G[H]$ in the branch; *a k -plex is in the branch (S, C) if and only if $S \subseteq H \subseteq S \cup C$* . To solve a branch (S, C) , it recursively solves two sub-branches formed based on a *branching vertex* v selected from C : one branch $(S \cup \{v\}, C \setminus \{v\})$ includes v to the partial solution S (which finds the largest k -plex containing v in (S, C)), and the other $(S, C \setminus \{v\})$ discards v from the candidate set C (which finds the largest k -plex excluding v in (S, C)). Clearly, solving two formed sub-branches solves branch (S, C) , and solving the branch (\emptyset, V) finds the largest k -plex in G .

Our kPEX adopts a similar framework in [11], which is summarized in Algorithm 1 and involves two stages. **Stage-I** first includes, in Stage-I.1, a heuristic method called KPHeuris for computing a large k -plex $G[S^*]$ (maintained globally as the largest k -plex seen so far), which will be used to narrow down the search space (Line 1), and a reduction method called CF-CTCP for reducing the input graph G by removing unpromising vertices/edges that will not appear in any k -plex larger than $|S^*|$ (Line 2). Besides, kPEX employs a widely-used *divide-and-conquer* strategy in Stage-I.2, which divides the problem of finding the largest k -plex in G into several sub-problems (Lines 3-7). Each sub-problem corresponds to a vertex v in G and aims to find the largest k -plex that (1) includes vertex v and (2) is in a subgraph of G induced by v 's two-hop neighbours $N^{\leq 2}(v)$, i.e., the set of vertices that have distance at most 2 from v (note that a k -plex with at least $2k - 1$ vertices has the diameter of at most 2 [63] and thus the largest k -plex containing v is a subset of $N^{\leq 2}(v)$). Clearly, the largest k -plex in G is the largest one among those returned by all sub-problems. **Stage-II** corresponds to the recursive process of solving a branch (Lines 9-15). Specifically, BRB_Rec recursively branches as discussed above (Lines 13-15). Besides, BRB_Rec conducts the newly proposed *alternated reduction-and-bound* process (AltRB) on a branch (S, C) for *narrowing down* the search space (Line 10). Specifically, it refines C to C^* by removing some unpromising vertices and computes an upper bound UB^* of (the size of) the largest k -plex in (S, C) for terminating the branch. Finally, we can terminate the branch when (1) $UB^* \leq |S^*|$ since no larger k -plex is in the branch and (2) $S \cup C^*$ is a k -plex since $G[S \cup C^*]$ is the largest k -plex in the branch.

Novelty. Our framework differs from the state-of-the-art one [11] in the following aspects. **First**, in Stage-II, kPEX is based on the newly proposed AltRB for narrowing down the search space. Recall that existing methods conduct the reduction-and-bound (RB) process using a sequential method called SeqRB at Line 10 instead. We will show that AltRB performs better than SeqRB in Section 4. Specifically, it refines C to a *smaller* set C^* (i.e., $|C^*| \leq |C|$) and obtains a *tighter* upper bound UB^* (i.e., $UB^* \leq UB$). **Second**, in Stage-I.1, kPEX employs the novel KPHeuris and CF-CTCP which are more effective and efficient than existing competitors in Section 5.

4 OUR REDUCTION&BOUND METHOD: ALTRB

4.1 An Alternated Reduction-and-Bound Method

Recall that existing algorithms conduct the reduction-and-bound (RB) step using the sequential method SeqRB on a branch $B = (S, C)$ for narrowing down the search space. Specifically, SeqRB has two sequential procedures: 1) the *reduction process* refines the candidate set C to C' based on $|S^*|$ (i.e., the lower bound of the branch), i.e., removing from C those vertices that cannot appear in a k -plex larger than $|S^*|$; and 2) the *bounding process* obtains the upper bound of the largest k -plex in the refined branch (S, C') , i.e., the upper bound of the branch denoted by $UB(S, C')$. In this paper, we propose a new alternated reduction-and-bound method, called AltRB, which is based on a binary partition of a branch $B = (S, C)$ as below.

$$S = S_L \cup S_R, \quad C = C_L \cup C_R. \quad (1)$$

Algorithm 2: Alternated reduction-and-bound: AltRB

Input: A graph $G = (V, E)$, a branch (S, C) and an integer k

Output: Refined candidate set C^* and upper bound UB^*

```
1  $S_L, S_R, C_L, C_R \leftarrow \text{Partition}(G, S, C, k)$ ;  
2  $UB_L \leftarrow |C_L|, LB_L \leftarrow 0$ ;  
3 while  $UB_L$  is not equal to  $\text{ComputeUB}(S_L, C_L)$  do  
4    $UB_L \leftarrow \text{ComputeUB}(S_L, C_L)$ ;  
5    $LB_R \leftarrow (|S^*| + 1) - |S| - UB_L$ ;  $C_R \leftarrow \text{RR1\&RR2}$  on  $C_R$ ;  
6    $UB_R \leftarrow \text{ComputeUB}(S_R, C_R)$ ;  
7    $LB_L \leftarrow (|S^*| + 1) - |S| - UB_R$ ;  $C_L \leftarrow \text{RR1\&RR2}$  on  $C_L$ ;  
8 return  $C^* \leftarrow C_L \cup C_R$  and  $UB^* \leftarrow |S| + UB_L + UB_R$ ;
```

Let $G[H]$ be a k -plex in the branch B such that $G[H]$ is larger than the largest k -plex $G[S^*]$ seen so far, i.e., $|H| \geq |S^*| + 1$ (note that other k -plexes have the size at most $|S^*|$ and thus can be ignored during the exploration of the branch). Based on the above partition, a k -plex $G[H]$ in B can be divided into three parts as below.

$$H = S \cup (C_L \cap H) \cup (C_R \cap H). \quad (2)$$

We denote by LB_L and UB_L (resp. LB_R and UB_R) the lower and upper bounds of the size of $C_L \cap H$ (resp. $C_R \cap H$), respectively. Formally, we have

$$|C_L \cap H| \leq UB_L, |C_R \cap H| \leq UB_R. \quad (3)$$

Besides, we have the following lemma on the above partition.

LEMMA 4.1. *Given a branch (S, C) with a partition, we have*

$$|C_L \cap H| \geq (|S^*| + 1) - |S| - UB_R, |C_R \cap H| \geq (|S^*| + 1) - |S| - UB_L. \quad (4)$$

PROOF. This can be easily verified since otherwise if $|C_L \cap H| < (|S^*| + 1) - |S| - UB_R$, we have $|H| = |S| + |C_L \cap H| + |C_R \cap H| < |S| + (|S^*| + 1) - |S| - UB_R + UB_R = |S^*| + 1$, which contradicts with $|H| \geq |S^*| + 1$. A similar contradiction can be derived for the other case $|C_R \cap H| < (|S^*| + 1) - |S| - UB_L$. \square

Based on Lemma 4.1, we define LB_L and LB_R as follows.

$$(|S^*| + 1) - |S| - UB_R \leq LB_L \leq |C_L \cap H| \quad (5)$$

$$(|S^*| + 1) - |S| - UB_L \leq LB_R \leq |C_R \cap H| \quad (6)$$

We note that Lemma 4.1 and Equations (5) and (6) indicate the relation between the lower bound of one part and the upper bound of the other, which enables AltRB. We summarize AltRB in Algorithm 2, which *iteratively and alternatively* conducts the reduction-and-bound step on the two partitions obtained via Partition (Line 1). Specifically, after initializing UB_L and LB_L in Line 2, AltRB involves the following steps (the details of the two procedures Partition and ComputeUB are provided in Section 4.2).

- **Step 1 (Bound on C_L).** Compute the upper bound for C_L (i.e., UB_L) via a procedure ComputeUB (Line 4).
- **Step 2 (Reduction on C_R).** Update the lower bound for C_R (i.e., LB_R) by $(|S^*| + 1) - |S| - UB_L$ according to Lemma 4.1 and then refine C_R based on the updated bounds via reduction rules RR1 and RR2 (Line 5).
- **Step 3 (Bound on C_R).** Compute the upper bound for the refined C_R (i.e., UB_R) via a procedure ComputeUB (Line 6).

- **Step 4 (Reduction on C_L).** Update the lower bound for C_L (i.e., LB_L) by $(|S^*| + 1) - |S| - UB_R$ according to Lemma 4.1 and then refine C_L based on the updated bounds via reduction rules RR1 and RR2 (Line 7).

Finally, we repeat Steps 1-4 until UB_L remains unchanged (Line 3). We remark that once tighter upper bounds are obtained at Step 1 and Step 3, tighter lower bounds can be derived via Lemma 4.1 at Step 2 and Step 4 which will be used to boost the performance of RR1 and RR2. Below find the details of reduction rules.

RR1. Given a branch (S, C) with LB_L and LB_R , 1) for a vertex v in C_L , we remove v from C if $|N(v, S \cup C_L)| < LB_L + |S| - k$ or $|N(v, S \cup C_R)| < LB_R + |S| - k + 1$; and 2) for a vertex v in C_R , we remove v from C if $|N(v, S \cup C_L)| < LB_L + |S| - k + 1$ or $|N(v, S \cup C_R)| < LB_R + |S| - k$.

RR2. Given a branch (S, C) with UB_L and UB_R , 1) if $UB_L + UB_R + |S| = |S^*| + 1$ and $UB_L = |C_L|$, we move all vertices in C_L from C to S if $G[S \cup C_L]$ is a k -plex; otherwise, i.e., it is not a k -plex, we terminate the branch (S, C) ; 2) if $UB_L + UB_R + |S| = |S^*| + 1$ and $UB_R = |C_R|$, we move all vertices in C_R from C to S if $G[S \cup C_R]$ is a k -plex; otherwise, i.e., it is not a k -plex, we terminate the branch (S, C) .

Benefits. Before proving the correctness, we show that AltRB better narrows down the search space than the existing SeqRB. The rationale behind is based on the following observations. First, at **Step 2** and **Step 4**, RR1, and RR2 (which are based on UB_L , UB_R , LB_L and LB_R) will remove from C more vertices when the lower bounds LB_L and LB_R become larger and/or the upper bounds UB_L and UB_R become smaller; Second, at **Step 1** and **Step 3**, with some vertices being removed from C_L and C_R , smaller upper bound UB_L and UB_R can be derived via ComputeUB (details refer to Section 4.2), and larger lower bounds LB_L and LB_R can also be obtained via Lemma 4.1; Third, as AltRB iteratively proceeds, the bounding process and the reduction process will benefit each other (since the former will derive smaller upper bounds and larger lower bounds after the latter while the latter will remove more vertices from C after the former). In contrast, SeqRB cannot be conducted iteratively since (1) its reduction rules are only based on $|S^*|$, which will not be changed after SeqRB and (2) thus repeating it multiple times cannot result in either a smaller candidate set C or a smaller upper bound. We remark that the refined set C^* and the upper bound UB^* obtained by AltRB is potentially *smaller* than those obtained by SeqRB (which will be proved in Section 4.2). Thus, with the proposed AltRB, our algorithm kPEX runs up to *two orders of magnitude faster* than the state-of-the-arts, as verified in our experiments.

Correctness. We then show the correctness of AltRB. Note that AltRB admits an arbitrary partition on (S, C) and any possible procedure for computing UB_L and UB_R that satisfy Equation (3).

The correctness of RR1 can be proved by contradiction. Consider a k -plex $G[H]$ in branch B with $|H| \geq |S^*| + 1$. Note that if such a k -plex does not exist, RR1 is obviously correct since all k -plexes in branch B are no larger than $|S^*|$ and thus branch B can be terminated. In general, there are two cases. First, assume that $G[H]$ contains a vertex v in C_L such that $|N(v, S \cup C_L)| < LB_L + |S| - k$. We get the contradiction by showing that v has more than k non-neighbours in H and thus $G[H]$ is not a k -plex since $|N(v, H)| = |N(v, H \cap (S \cup C_L))| + |N(v, H \cap C_R)| \leq (LB_L + |S| - k - 1) + |H \cap C_R| \leq$

$(|S| + |H \cap C_R| + |H \cap C_L|) - (k+1) = |H| - (k+1)$. **Second**, assume that $G[H]$ contains a vertex v in C_L such that $|N(v, S \cup C_R)| < LB_R + |S| - k + 1$. Similarly, we derive the contradiction by showing that v has more than k non-neighbours in H and thus $G[H]$ is not a k -plex since $|N(v, H)| = |N(v, H \cap (S \cup C_R))| + |N(v, H \cap C_L)| \leq (LB_R + |S| - k) + (|H \cap C_L| - 1) \leq (|S| + |H \cap C_R| + |H \cap C_L|) - (k+1) = |H| - (k+1)$ (note that $|N(v, H \cap C_L)| \leq |H \cap C_L| - 1$ since v is in C_L and is not adjacent to itself). Symmetrically, we can prove the correctness for the reduction rules on C_R .

The correctness of **RR2** is easy to verify. Consider a branch (S, C) with $UB_L + UB_R + |S| = |S^*| + 1$ and $UB_L = |C_L|$, and a k -plex $G[H]$ in (S, C) with $|H| \geq |S^*| + 1$ (note that if such a k -plex does not exist, **RR2** is obviously correct on this branch). We note that $G[H]$ must contain all vertices in C_L , i.e., $C_L \subseteq H$, since otherwise $|H| = |H \cap S| + |H \cap C_L| + |H \cap C_R| \leq |S| + (|C_L| - 1) + |H \cap C_R| \leq |S| + UB_L + UB_R - 1 = |S^*|$. Therefore, $G[S \cup C_L]$ must be a k -plex due to the hereditary property; otherwise, such a k -plex cannot exist in (S, C) and we can terminate the branch.

The correctness of **AltRB** can then be easily verified.

4.2 Upper Bound Computation and Greedy Partition Strategy

In this part, we first introduce the method `ComputeUB` used at **Step 1** and **Step 3** for obtaining UB_L and UB_R in Section 4.1. To boost the performance of `ComputeUB` as well as the reduction rules on C_L and C_R , we then propose a greedy strategy `Partition` for partitioning C (resp. S) into C_L and C_R (resp. S_L and S_R). Finally, with all carefully-designed techniques above, we show that the resulted upper bound UB^* will be potentially *smaller* than the existing one UB .

Upper bound computation. We adapt an existing upper bound computation [35], which we call `ComputeUB`, for obtaining UB_L and UB_R . Note that it can handle an arbitrary partition on a branch (S, C) . Consider **Step 1** for computing UB_L . `ComputeUB`(S_L, C_L) first iteratively partitions C_L into $(|S_L| + 1)$ disjoint subsets. The i -th ($1 \leq i \leq |S_L|$) subset $\Pi_i(S_L, C_L)$ contains all non-neighbours of a vertex $u_i \in S_L$ in $C_L - \{\Pi_1(S_L, C_L), \dots, \Pi_{i-1}(S_L, C_L)\}$, formally,

$$\Pi_i(S_L, C_L) = \overline{N}(u_i, C_L^i), \quad C_L^i = C_L - \bigcup_{j=1}^{i-1} \Pi_j(S_L, C_L), \quad (7)$$

where u_i is the vertex in $S_L \setminus \{u_1, u_2, \dots, u_{i-1}\}$ with the largest ratio of $|\overline{N}(u_i, C_L^i)| / (k - |\overline{N}(u_i, S)|)$. Note that the strategy of selecting u_i from S_L has been shown to boost the practical performance of `ComputeUB` (details refer to [35]). Besides, we have $\Pi_0(S_L, C_L) = C_L - \{\Pi_1(S_L, C_L), \dots, \Pi_{|S_L|}(S_L, C_L)\}$. Thus, vertices in $\Pi_i(S_L, C_L)$ ($1 \leq i \leq |S_L|$) are the non-neighbours of u_i in C_L , and vertices in $\Pi_0(S_L, C_L)$ are common neighbours of vertices in S_L . The key observation is that for a k -plex $G[H]$ in the branch, $C_L \cap H$ contains at most $\min\{|\Pi_i(S_L, C_L)|, k - |\overline{N}(u_i, S)|\}$ vertices from $\Pi_i(S_L, C_L)$ for $1 \leq i \leq |S_L|$ since otherwise u_i (in H) will have more than k non-neighbours in $G[H]$ and thus $G[H]$ is not a k -plex. Thus, the upper bound UB_L returned by `ComputeUB`(S_L, C_L) gives as below:

$$|\Pi_0(S_L, C_L)| + \sum_{i=1}^{|S_L|} \min\{|\Pi_i(S_L, C_L)|, k - |\overline{N}(u_i, S)|\}. \quad (8)$$

We note that with some vertices being removed from C_L during **AltRB**, $\Pi_i(S_L, C_L)$ will get smaller and thus a smaller upper bound can be derived. Similarly, we can obtain UB_R by `ComputeUB`(S_R, C_R).

Besides, we remark that the state-of-the-art upper bound of k -plex in the branch (S, C) (used in **SeqRB**) is $|S| + \text{ComputeUB}(S, C)$ [35].

Greedy partition. Consider the upper bound computation at C_L , i.e., Equation (8). We observe that *all vertices in $\Pi_0(S_L, C_L)$ contributes to the upper bound $\text{ComputeUB}(S_L, C_L)$* since each of them is adjacent to all vertices in S_L and thus they could appear in a k -plex in branch (S, C) . The similar observation can be derived on other subsets $\Pi_i(S_L, C_L)$ such that $|\overline{N}(u_i, C_L^i)| \leq k - |\overline{N}(u_i, S)|$ and $1 \leq i \leq |S_L|$ (note that there are *fewer* missing edges between S_L and those subsets). Therefore, the adapted upper bound computation performs worse on those subsets.

Motivated by the above observation, we propose to divide S and C into the one (S_L and C_L) with *more* missing edges and the other (S_R and C_R) with *fewer* missing edges. We summarize the proposed strategy in Algorithm 3. Specifically, we iteratively remove from S to S_L (resp. from C to C_L) the vertex v with the greatest value of $|\overline{N}(v, C)| / (k - |\overline{N}(v, S)|)$ (resp. the set of v 's non-neighbours in C , i.e., $\overline{N}(v, C)$) until the greatest value of $|\overline{N}(v, C)| / (k - |\overline{N}(v, S)|)$ is not greater than 1 or S becomes empty (Lines 2-6). Then, all remaining vertices in S and C will be removed to S_R and C_R (Line 7). We observe that (1) `ComputeUB`(S_L, C_L) will return a tighter bound since $|\overline{N}(u_i, C_L^i)| > k - |\overline{N}(u_i, S)|$ holds for $1 \leq i \leq |S_L|$ and $\Pi_0(S_L, C_L) = \emptyset$, and (2) `ComputeUB`(S_R, C_R) is always equal to $|C_R|$.

Consider a branch (S, C) (which has been refined by **SeqRB**) with the upper bound $UB = |S| + \text{ComputeUB}(S, C)$. With the proposed techniques, **AltRB** will further narrow down the search space of (S, C) by the following observation.

$$UB^* \leq UB \text{ and } |C^*| \leq |C|. \quad (9)$$

We note that $|C^*| \leq |C|$ is obvious since some vertices in C could be removed via **RR1** and **RR2**. Besides, $UB^* \leq UB$ holds since (1) `ComputeUB`(S, C) = `ComputeUB`(S_L, C_L) + `ComputeUB`(S_R, C_R) before **AltRB** (which can be verified based on the definitions) and (2) as **AltRB** proceeds, C_L and C_R are refined via **RR1** and **RR2**, and thus `ComputeUB`(S_L, C_L) and `ComputeUB`(S_R, C_R) get smaller.

Benefits of greedy partition. Compared with a random partition, the greedy partition in Algorithm 3 has the following advantageous properties. (1) A tight upper bound of C_L leads to a larger LB_R , which enhances the effectiveness of **RR1**. (2) $UB_R = |C_R|$ is always satisfied, which means that **RR2** is applicable as long as $UB_L + UB_R + |S| = |S^*| + 1$. In other words, the conditions for **RR2** are more relaxed. Moreover, computing UB_R as $|C_R|$ is easy to implement and requires less computation.

4.3 Time Complexity Analysis

Time complexity. We analyze the time complexity of **AltRB** as follows. (1) **AltRB** first invokes `Partition` (Algorithm 3). Specifically, Lines 2-6 of Algorithm 3 will be conducted at most $|S|$ times, and each iteration needs to compute $|\overline{N}(v, C)|$ for each $v \in S$, which can be done in $O(|S| \times |C|)$. Thus, the time complexity of `Partition` is $O(|S|^2 |C|)$. (2) **AltRB** then iteratively processes Lines 3-7 of Algorithm 2. We note that `ComputeUB`(S, C) can be computed in $O(|S|^2 |C|)$ [35] (Lines 4 and 6). For reductions rules in Lines 5 and 7, **RR1** iteratively removes the vertex in C_R with minimum $|N(v, S \cup C_L)|$ (or $|N(v, S \cup C_R)|$), and **RR2** checks whether $S \cup C_R$ is a k -plex. Both rules can be done in $O(|C| \times (|S| + |C|))$. (3) We

Algorithm 3: Partition(G, S, C, k)

Input: Branch (S, C) , a graph $G = (V, E)$, and an integer k

Output: The greedy partition S_L, S_R, C_L and C_R

```

1  $S_L \leftarrow \emptyset, S_R \leftarrow \emptyset, C_L \leftarrow \emptyset, C_R \leftarrow \emptyset;$ 
2 while  $S \neq \emptyset$  do
3    $v^* \leftarrow \arg \max_{v \in S} |\bar{N}(v, C)| / (k - |\bar{N}(v, S)|);$ 
4   if  $|\bar{N}(v^*, C)| / (k - |\bar{N}(v^*, S)|) \leq 1$  then break;
5    $S_L \leftarrow S_L \cup \{v^*\}, C_L \leftarrow C_L \cup \bar{N}(v^*, C);$ 
6    $S \leftarrow S \setminus \{v^*\}, C \leftarrow C \setminus \bar{N}(v^*, C);$ 
7  $S_R \leftarrow S, C_R \leftarrow C;$ 
8 return  $S_L, S_R, C_L$  and  $C_R;$ 

```

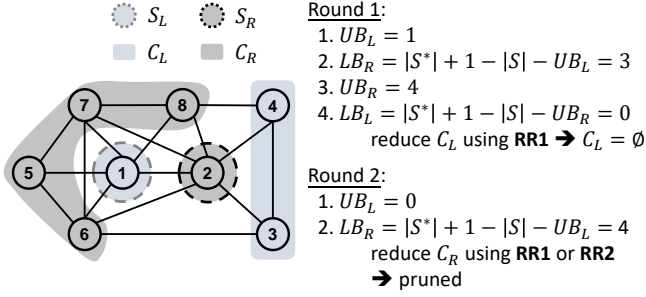


Figure 1: An example of AltrB with $k=2$, $|S^*|=5$, $S = \{v_1, v_2\}$, $C = \{v_3, v_4, v_5, v_6, v_7, v_8\}$

also know that $|S|$ is bounded by $\delta(G) + k$; otherwise, we have a k -plex $G[S]$ with $|S| > \delta(G) + k$, which will form a $(\delta(G) + 1)$ -core and thus contradict the definition of $\delta(G)$; $|C|$ is bounded by $\delta(G)d$ since $V(g)$ at Line 5 of Algorithm 1 is bounded by $\delta(G)d$ [17, 54]. (4) Let r be the number of iterations of Lines 3-7 of Algorithm 2. The number of r is quite small in practice (e.g., $r = 1.13$ on average in our experiments) and is bounded by $|C|$ (i.e., $\delta(G)d$) since at least one vertex is removed from C in each round until C becomes empty.

Thus, AltrB (Algorithm 2) runs in $O(r \times (|S|^2|C| + |C|^2)) = O(\delta(G)^3 d^3 + k^2 \delta(G)^2 d^2)$, where $\delta(G)$ is much smaller than d and n in real graphs, as shown in Table 1 ($\delta(G) \leq d < n$ in theory).

Example. To illustrate the proposed AltrB, consider an example in Figure 1 with $k=2$, $|S^*| = 5$, $S = \{v_1, v_2\}$ and $C = \{v_3, v_4, v_5, v_6, v_7, v_8\}$. First, we apply the greedy partition (Algorithm 3) and obtain $S_L = \{v_1\}$, $S_R = \{v_2\}$, $C_L = \{v_3, v_4\}$, and $C_R = \{v_5, v_6, v_7, v_8\}$. Then, in the first round of AltrB (Lines 4-7 in Algorithm 2), we conduct the four steps. (**Step 1**) Compute the upper bound of C_L , i.e., $UB_L = 1$. (**Step 2**) Update the lower bound of C_R (i.e., $LB_R = 3$) and reduce C_R via **RR1** and **RR2** (no vertices are removed). (**Step 3**) Compute the upper bound of C_R , i.e., $UB_R = 4$. (**Step 4**) Update the lower bound of C_L (i.e., $LB_L = 0$) and reduce C_L via **RR1** and **RR2** (C_L is reduced to an empty set). Next, in the second round with $C_L = \emptyset$, we (1) compute $UB_L = 0$, and (2) update $LB_R = 4$ and reduce C_R . If we first apply **RR1** to C_R , v_5, v_6, v_7 and v_8 will be removed, and finally compute the upper bound as $UB^* = |S| + |UB_L| + |UB_R| = 2 + 0 + 0 = 2$, resulting in pruning. If we first apply **RR2**, both $UB_L + UB_R + |S| = |S^*| + 1$ and $UB_R = |C_R|$ are satisfied. We then find that $G[S \cup C_R]$ is not

a k -plex, which means that **RR2** also leads to pruning. Actually, the size of maximum 2-plex is 5, indicating that the branch (S, C) cannot find a larger 2-plex, and thus this branch can be pruned by AltrB. However, without AltrB, the existing method [35] will compute an upper bound as $UB = UB_L + UB_R + |S| = 1 + 4 + 2 = 7$, which cannot prune the current branch.

Remarks. We remark that the existing reduction rules proposed in [11, 12, 53] are all based on $|S^*|$ and thus orthogonal to AltrB. We conduct some of these reduction rules to improve practical performance, including (1) additional reduction on subgraph g (Lemma 3.2 in [11] and Reduction 2 in [53]) in Line 5 of Algorithm 1, and (2) reduction on C before AltrB (**RR4** in [11] and Algorithm 3 in [12]). Besides, AltrB is also orthogonal to the branching rules for selecting the branching vertex and forming the sub-branches.

5 EFFICIENT PRE-PROCESSING TECHNIQUES

In this section, we develop some efficient pre-processing techniques for further boosting the performance of BRB algorithms, namely, CF-CTCP for reducing the size of the input graph in Section 5.1 and KPHeur is for heuristically computing a large k -plex in Section 5.2.

5.1 Faster Core-Truss Co-Pruning: CF-CTCP

Let lb be the lower bound of the size of the largest k -plex (which corresponds to the size of the largest k -plex $G[S^*]$ seen so far). We also let $\Delta(u, v)$ be the set of common neighbors of u and v in G , i.e., $\Delta(u, v) = N_G(u) \cap N_G(v)$. The idea of refining the input graph G is to remove from G those vertices and edges that cannot appear in any k -plex larger than lb as many as we can. Existing methods [11, 35, 63] are all based on the following lemmas and differ in the implementations (the details of proof is omitted for the ease of presentation).

LEMMA 5.1. (Core Pruning [27]) For each vertex $u \in V(G)$, u cannot appear in a k -plex of size $lb + 1$ if $d_G(u) \leq lb - k$.

LEMMA 5.2. (Truss Pruning [63]) For each edge $(u, v) \in E(G)$, (u, v) cannot appear in a k -plex of size $lb + 1$ if $\delta_G(u, v) \leq lb - 2k$ where $\delta_G(u, v)$ is the number of common neighbors of u and v , i.e., $\delta_G(u, v) = |\Delta(u, v)|$.

Note that the time complexities of core pruning and truss pruning are $O(m)$ [4] and $O(m \times \delta(G))$ [51], respectively. The above lemmas (namely, core pruning and truss pruning) indicates those unpromising vertices and edges can be removed from G . In particular, with some vertices or edges being removed from G , the remaining vertices u and edges (u, v) have $d_G(u)$ and $\delta_G(u, v)$ decreases, respectively; and then more vertices and edges can be removed. Therefore, the core pruning (resp. the truss pruning) can be conducted in an iterative way, i.e., iteratively removing unpromising vertices (resp. edges) and updating $d_G(\cdot)$ (resp. $\delta_G(\cdot, \cdot)$) for the remaining until no vertex or edge can be removed. We remark that the state-of-the-art method called the core-truss co-pruning (CTCP [11]) iteratively conducts the truss pruning and then the core pruning in multiple rounds until the graph remains unchanged. However, we observe that CTCP is still inefficient due to potential redundant computations. This is because (1) CTCP performs the truss pruning and the core pruning *separately* at each round (i.e., first remove a set of edges via the truss pruning and then remove

one unpromising vertex via core pruning), (2) the truss pruning has the time complexity of $O(m \times \delta(G))$ larger than $O(m)$ for the core pruning, and (3) we note that during the truss pruning, some vertices can be removed via the more efficient core pruning while the truss pruning will iteratively check all their incident edges and then remove some of them (which is very costly).

To improve the practical efficiency of CTCP, we propose a new algorithm called the *core-pruning-first core-truss co-pruning* (or CF-CTCP), which differs from CTCP in the way of conducting pruning at each round. Specifically, at each round, it first removes *all* unpromising vertices and then removes *one* unpromising edge (recall that CTCP first removes *all* unpromising edges and then *one* unpromising vertex). The benefit is that unpromising vertices can be removed immediately via efficient core pruning. Note that our CF-CTCP has the same output but requires less computation compared to CTCP. Given the integer k and the lower bound lb , both CF-CTCP and CTCP reduce the input graph G to the maximal subgraph that is $(lb+1-k)$ -core and $(lb+3-2k)$ -truss. The difference between CF-CTCP and CTCP is illustrated in Figure 2.

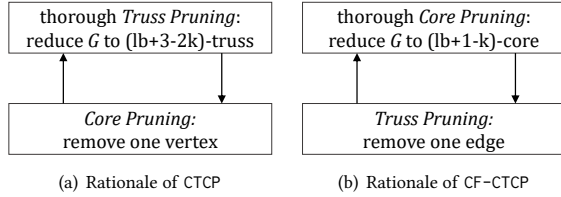


Figure 2: Comparing CTCP and CF-CTCP

The main idea of CF-CTCP is to conduct core pruning thoroughly as follows: 1) if we identify an edge that can be removed, we will immediately remove this edge, even if we have not yet finished computing $\Delta(\cdot, \cdot)$ (i.e., all triangles for each edge); 2) after removing an edge (u, v) , we will check whether u or v can be reduced by core pruning. Note that after removing an edge (u, v) , we postpone the action of updating $\Delta(u, \cdot)$ and $\Delta(v, \cdot)$ since it is time-consuming and there may lead to redundant computations. For example, if both vertices u and v will be removed by core pruning later, updating $\Delta(u, \cdot)$ and $\Delta(v, \cdot)$ is not necessary.

Our proposed CF-CTCP is shown in Algorithm 4. The input of CF-CTCP includes: 1) a set of vertices Q_v , which stores the vertices that need to be removed; 2) two integers $\tau_v = lb - k$ and $\tau_e = lb - 2k$ that serve as thresholds for the numbers of degrees and triangles for pruning, respectively; 3) a boolean value $lb_changed$ which is *true* if a larger k -plex is found in kPEX and KPHeur is (Algorithm 1 and Algorithm 5). We note that both kPEX and KPHeur is (Algorithms 1 and 5) invoke CF-CTCP multiple times. For example, KPHeur is invokes CF-CTCP by calling $CF-CTCP(G, \emptyset, lb - k, lb - 2k, true)$ when it finds a larger heuristic k -plex of size lb .

We then describe the details of CF-CTCP in steps. First, we design a procedure called *RemoveEdge* (Lines 21-24) to remove one unpromising edge in Line 21 and all current unpromising vertices in Line 22 via core and truss pruning. The set of removed edges to be considered (due to Lines 21 and 22) is pushed into Q_e , which will be used to update $\Delta(\cdot, \cdot)$ later. Second, Lines 5-6 initialize the

Algorithm 4: CF-CTCP($G = (V, E), Q_v, \tau_v, \tau_e, lb_changed$)

Input: A graph $G = (V, E)$, the set of vertices to be removed Q_v , two integral thresholds τ_v and τ_e , a boolean value $lb_changed$

Output: The reduced graph which is the maximal subgraph in G that is both a $(\tau_v + 1)$ -core and a $(\tau_e + 3)$ -truss

```

1 Remove the vertices in  $Q_v$  from  $G$  and reduce  $G$  to the
  maximal  $(\tau_v + 1)$ -core by the core pruning;
2 Initialize the set of removed edges to be considered
   $Q_e \leftarrow \{\text{edges removed at Line 1}\};$ 
3 if  $lb\_changed$  then
4   for each  $(u, v) \in E$  do
5     if CF-CTCP is invoked for the first time then
6        $\Delta(u, v) \leftarrow N_G(u) \cap N_G(v);$ 
7     if  $|\Delta(u, v)| \leq \tau_e$  then
8        $Q_e \leftarrow Q_e \cup \text{RemoveEdge}(G, (u, v), \tau_v);$ 
9 while  $Q_e \neq \emptyset$  do
10    $(u, v) \leftarrow \text{pop an edge from } Q_e;$ 
11   if  $u \in V$  then
12     for each  $w \in N_G(u)$  satisfying  $v \in \Delta(u, w)$  do
13       Remove  $v$  from  $\Delta(u, w);$ 
14       if  $|\Delta(u, w)| \leq \tau_e$  then
15          $Q_e \leftarrow Q_e \cup \text{RemoveEdge}(G, (u, w), \tau_v);$ 
16   if  $v \in V$  then
17     for each  $w \in N_G(v)$  satisfying  $u \in \Delta(v, w)$  do
18       Remove  $u$  from  $\Delta(v, w);$ 
19       if  $|\Delta(v, w)| \leq \tau_e$  then
20          $Q_e \leftarrow Q_e \cup \text{RemoveEdge}(G, (v, w), \tau_v);$ 

```

Procedure: *RemoveEdge*($G, (u, v), \tau_v$)

Output: The set of removed edges to be considered Q_e

```

21 Remove the unpromising edge  $(u, v)$  from  $G;$ 
22 Reduce  $G$  to the maximal  $(\tau_v + 1)$ -core by the core pruning;
23 Initialize the set of removed edges to be considered
   $Q_e \leftarrow \{\text{edges removed at Lines 21-22}\};$ 
24 return  $Q_e;$ 

```

sets of common neighbours $\Delta(\cdot, \cdot)$ if CF-CTCP is invoked for the first time. Whenever we find an edge (u, v) that can be reduced, we invoke the procedure *RemoveEdge* to remove (u, v) immediately in Line 8. Third, we postpone the action of updating $\Delta(\cdot, \cdot)$ to Lines 9-20. Lines 11-20 consider the effect of each removed edge (u, v) by traversing all the triangles that (u, v) participates in. Specifically, Lines 11-15 traverse each edge $(u, w) \in E$ satisfying $v \in \Delta(u, w)$, i.e., u, v, w can form a triangle, then we update $\Delta(u, w)$ and check whether edge (u, w) can be reduced. Lines 16-20 consider the edges connected to v , which is similar to Lines 11-15. Note that in Lines 15 and 20, if we find an edge that can be reduced, we invoke the procedure *RemoveEdge* to remove the edge immediately.

Time complexity. We analyze the time complexity of CF-CTCP (Algorithm 4), including all invocations in kPEX, in the following.

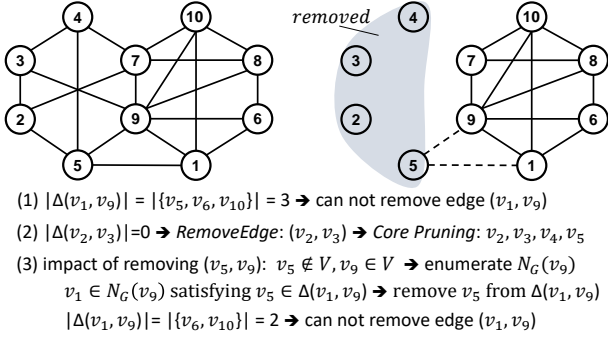


Figure 3: An example for CF-CTCP assuming $lb = 4$ and $k = 2$

LEMMA 5.3. *The total time complexity of all invocations in kPEX (Algorithm 1 which includes invocations in the heuristic process KPHeur is in Algorithm 5) to CF-CTCP (Algorithm 4) is $O(m \times \delta(G))$.*

The omitted proof, along with an implementation of CF-CTCP with $O(m)$ memory usage, is provided in the appendix of the technical report [29].

Remarks. First, the time complexity of CTCP is $O(m \times \delta(G) + m \times k) = O(m\delta(G))$, requiring that k is a small constant. However, k is up to n in theory and the time complexity of our CF-CTCP is always $O(m\delta(G))$ for all possible values of k . **Second**, we do not consider the update of $\Delta(\cdot, \cdot)$ when removing a vertex because removing a vertex is equivalent to first removing all the edges connected to this vertex and then removing this isolated vertex. Therefore, we only consider the removed edges for updating $\Delta(\cdot, \cdot)$. **Third**, the acceleration of CF-CTCP can be attributed to two main factors: 1) we do not need to compute the numbers of triangles for the edges that can be removed by core pruning; 2) for an edge (u, v) to be removed such that both u and v are already removed by core pruning, we do not need to traverse related triangles to update $\Delta(\cdot, \cdot)$. Note that if we cannot remove any vertex or edge, the time consumption of CF-CTCP will be the same as CTCP in theory, which is due to the fact that both of them need to compute $\Delta(\cdot, \cdot)$ in $O(m \times \delta(G))$.

An example of CF-CTCP. Consider the example of CF-CTCP (Algorithm 4) in Figure 3, assuming $lb = 4$ and $k = 2$. According to Lemma 5.1 and Lemma 5.2, we need to reduce G to the maximal subgraph that is both a 3-core and a 3-truss, i.e., we will remove a vertex u if $d_G(u) < 3$ and an edge (u, v) if $|\Delta(u, v)| < 1$. **First**, we enumerate each edge (u, v) and compute the common neighbors of u and v (Lines 4-6). For those edges connected to v_1 , we cannot remove them since they have enough common neighbors, e.g., there are 3 common neighbors of v_1 and v_9 . However, when we consider edges connected to v_2 , we find that edge (v_2, v_3) can be removed since $|\Delta(v_2, v_3)| = 0$. We then immediately remove edge (v_2, v_3) and conduct core pruning, which removes vertices v_2, v_3, v_4 , and v_5 (Lines 21-24). After this process, we continue to compute common neighbors for the remaining edges, but none of these edges can be removed. **Second**, we begin to consider those removed edges in Q_e . We focus on the edges (v_5, v_9) and (v_1, v_5) since the other removed edges cannot form a triangle with the remaining edges in G . For the removal of the edge (v_5, v_9) , according to Lines 11-20, we update $\Delta(v_1, v_9)$, and the triangle (v_1, v_5, v_9) is destroyed. Then,

for the edge (v_1, v_5) , since the triangle (v_1, v_5, v_9) no longer exists after removing the edge (v_5, v_9) , the edge (v_1, v_5) cannot constitute any triangle with other vertices. Thus, the procedure of CF-CTCP terminates. Finally, we reduce G to $G[\{v_1, v_6, v_7, v_8, v_9, v_{10}\}]$ where $G[\{v_1, v_6, v_8, v_9, v_{10}\}]$ is a 2-plex of size 5.

5.2 Compute a large k -plex: KPHeur is

We introduce a heuristic method KPHeur is for computing a large initial k -plex. Note that such an initial k -plex offers a lower bound lb , which helps to narrow the search space; and the larger the lower bound is, the more search space can be refined. Therefore, KPHeur is is designed for obtaining a large k -plex *efficiently and effectively*.

Algorithm 5: KPHeur is(G, k)

Input: A graph $G = (V, E)$ and an integer $k > 1$

Output: The vertex set S of a heuristic initial k -plex $G[S]$

```

1  $S \leftarrow \text{Degen}(G, k, lb \leftarrow |S|)$ ;
2 Apply CF-CTCP for refining  $G$  based on  $lb$ ;
3 for each  $v_i \in V(G)$  do
4    $g \leftarrow G[\{v_i, v_{i+1}, \dots, v_n\} \cap N^{\leq 2}(v_i)]$ ;  $S' \leftarrow \text{Degen}(g, k)$ ;
5   if  $|S'| > |S|$  then
6      $S \leftarrow S', lb \leftarrow |S|$ ;
7   Apply CF-CTCP for refining  $G$  based on  $lb$ ;
8 return  $S$ ;

Procedure: Degen( $G, k$ )
Output: The vertex set  $S$  of a heuristic maximal  $k$ -plex in  $G$ 
9  $v_1, v_2, \dots, v_n \leftarrow$  the degeneracy order of vertices in  $V(G)$ ;
10  $S \leftarrow \emptyset$ ;
11 for  $i = n$  to 1 do
12   if  $G[S \cup \{v_i\}]$  is a  $k$ -plex then  $S \leftarrow S \cup \{v_i\}$ ;
13 return  $S$ ;
```

We summarize KPHeur is in Algorithm 5, which relies on a sub-procedure (called Degen) for computing a large k -plex. Specifically, Degen iteratively includes to an empty set S a vertex in a graph G based on the degeneracy ordering while retaining the k -plex property of $G[S]$ until we cannot continue (Lines 9-12). To compute a larger k -plexes, KPHeur is further generate n subgraphs from G , each of which corresponds to a vertex in G (Lines 3-4); it then invokes Degen on each of them to obtain a k -plex (Line 4); it finally returns the largest one among $n + 1$ found k -plexes. Note that the subgraph related to v_i is the subgraph induced by $\{v_i, v_{i+1}, \dots, v_n\} \cap N^{\leq 2}(v_i)$ where $N^{\leq 2}(u)$ denotes u 's neighbors and u 's neighbors' neighbors, and the rationale is that it can make the subgraph smaller and denser, which tends to find a larger k -plex easier. The time complexity of Degen is $O(m)$, and we will invoke it at most $n + 1$ times, thus the total time complexity of computing heuristic solutions in Algorithm 5 is $O(nm)$. We remark that the total time complexity of all invocations of CF-CTCP is $O(m\delta(G))$ because we invoke CF-CTCP in KPHeur is only when we find a larger k -plex, i.e., $lb_changed = \text{true}$, as shown in Lemma 5.3. Thus, the time complexity of KPHeur is is $O(m\delta(G) + nm) = O(nm)$.

Compared with existing heuristic methods. There are two state-of-the-art heuristic methods: kPlex-Degen ([11]) and EGo-Degen

([12]). kPlex-Degen computes a large k -plex by iteratively removing a vertex from the input graph G based on a certain ordering until the remaining graph becomes a k -plex. KPHeur is differs from kPlex-Degen in two aspects. First, Degen computes a large k -plex by iteratively including a vertex, which is more efficient since the size of the largest k -plex is usually much smaller than the size of the input graph (especially for real-world graphs) and can always return a maximal k -plex, while kPlex-Degen cannot. Second, KPHeur is further explores n subgraphs instead of the input graph G , which tends to obtain a larger k -plex as empirically verified in our experiments. EGo-Degen extracts a subgraph g_v for each vertex v and invokes kPlex-Degen to compute a k -plex in g_v . Then, EGo-Degen selects the largest k -plex among those computed on n subgraphs as the initial heuristic k -plex. KPHeur is differs from EGo-Degen in three aspects. First, the method of subgraph extraction is different. For a vertex $v \in V(G)$, EGo-Degen extracts $g_v = G[\{v_i, v_{i+1}, \dots, v_n\} \cap N_G(v_i)]$, while our KPHeur is generates a subgraph $g'_v = G[\{v_i, v_{i+1}, \dots, v_n\} \cap N^{\leq 2}(v_i)]$. It is easy to verify that $g_v \subseteq g'_v$ due to $N_G(v_i) \subseteq N^{\leq 2}(v_i)$. Additionally, a larger subgraph tends to contain a larger k -plex, as verified in Table 6. Second, EGo-Degen computes k -plexes by invoking kPlex-Degen, which implies that it may find a non-maximal k -plex as mentioned above. Third, once a larger k -plex is found, KPHeur is updates lb and removes unpromising vertices/edges immediately, while EGo-Degen does not reduce the graph until n heuristic k -plexes are computed.

6 EXPERIMENTAL STUDIES

We test the efficiency and effectiveness of our algorithm kPEX by comparing with the state-of-the-art BRB algorithms:

- **kPlexS**¹: the existing algorithm proposed in [11].
- **KPLEX**²: the existing algorithm proposed in [53].
- **DiseMKP**³: the existing algorithm proposed in [35].
- **kPlexT**⁴: the existing algorithm proposed in [12].

Setup. All algorithms are written in C++ and compiled with -O3 optimization by g++ 9.4.0. Moreover, all algorithms are initialized with a lower bound of $2k - 2$ to focus on finding k -plexes with at least $2k - 1$ vertices. All experiments are conducted in the single-thread mode on a machine with an Intel(R) Xeon(R) Platinum 8358P CPU@2.60GHz and 256GB main memory. The CPU frequency is fixed at 3.3GHz. We set the time limit as 3600 seconds and use **OOT** (Out Of Time limit) to indicate the time exceeds the limit. We consider six different values of k , i.e., 2, 3, 5, 10, 15, and 20. We focus on the case of $k = 5$, and defer the experiments for other values of k to the appendix of the technical report [29]. We also note that the major findings for $k = 5$ hold for other values of k .

Datasets. We consider the following two collections of graphs.

- **Network Repository** [2]. The dataset contains 584 graphs with up to 5.87×10^7 vertices, including biological networks (36), dynamic networks (85), labeled networks (104), road networks (15), interaction (29), scientific computing (11), social networks (75), facebook (114), web (31), and DIMACS-10 graphs (84). Most of them are real-world graphs.

Table 1: Statistics of 30 representative graphs

ID	Graph	n	m	density	d_{max}	$\delta(G)$
G1	johnson8-4-4	70	1855	$7.68 \cdot 10^{-1}$	53	53
G2	C125-9	125	6963	$8.98 \cdot 10^{-1}$	119	102
G3	keller4	171	9435	$6.49 \cdot 10^{-1}$	124	102
G4	brock200-2	200	9876	$4.96 \cdot 10^{-1}$	114	84
G5	san200-0-9-1	200	17910	$9.00 \cdot 10^{-1}$	191	162
G6	san200-0-9-2	200	17910	$9.00 \cdot 10^{-1}$	188	169
G7	san200-0-9-3	200	17910	$9.00 \cdot 10^{-1}$	187	169
G8	p-hat300-1	300	10933	$2.44 \cdot 10^{-1}$	132	49
G9	p-hat300-2	300	21928	$4.89 \cdot 10^{-1}$	229	98
G10	p-hat500-1	500	31569	$2.53 \cdot 10^{-1}$	204	86
G11	soc-BlogCatalog-ASU	10312	333983	$6.28 \cdot 10^{-3}$	3992	114
G12	socfb-Ullinois	30795	1264421	$2.67 \cdot 10^{-3}$	4632	85
G13	soc-themarker	69413	1644843	$6.83 \cdot 10^{-4}$	8930	164
G14	soc-BlogCatalog	88784	2093195	$5.31 \cdot 10^{-4}$	9444	221
G15	soc-buzznet	101163	2763066	$5.40 \cdot 10^{-4}$	64289	153
G16	soc-LiveMocha	104103	2193083	$4.05 \cdot 10^{-4}$	2980	92
G17	soc-wiki-conflict	116836	2027871	$2.97 \cdot 10^{-4}$	20153	145
G18	soc-google-plus	211187	1141650	$5.12 \cdot 10^{-5}$	1790	135
G19	soc-FourSquare	639014	3214986	$1.57 \cdot 10^{-5}$	106218	63
G20	rec-epinions-user-ratings	755760	13667951	$4.79 \cdot 10^{-5}$	162179	151
G21	soc-wiki-Talk-dir	1298165	2288646	$2.72 \cdot 10^{-6}$	100025	119
G22	soc-pokec	1632803	22301964	$1.67 \cdot 10^{-5}$	14854	47
G23	tech-ip	2250498	21643497	$8.55 \cdot 10^{-6}$	1833161	253
G24	ia-wiki-Talk-dir	2394385	4659565	$1.63 \cdot 10^{-6}$	100029	131
G25	sx-stackoverflow	2584164	28183518	$8.44 \cdot 10^{-6}$	44065	198
G26	web-wikipedia_link_it	2790239	86754664	$2.23 \cdot 10^{-5}$	825147	894
G27	socfb-A-anon	3097165	23667394	$4.93 \cdot 10^{-6}$	4915	74
G28	soc-livejournal-user-groups	7489073	112305407	$4.00 \cdot 10^{-6}$	1053720	116
G29	soc-bitcoin	24575382	86063840	$2.85 \cdot 10^{-7}$	1083703	325
G30	soc-sinaweibo	58655849	261321033	$1.52 \cdot 10^{-7}$	278489	193

- **2nd-DIMACS (DIMACS-2) Graphs** [1]. The dataset contains 80 synthetic dense graphs with up to 4000 vertices and the densities ranging from 0.03 to 0.99. Most graphs in the dataset are synthetic graphs, which are often hard to be solved [35, 36, 53].

For better comparisons, we select 30 representative graphs from the above 664 graphs and report the statistics in Table 1, where the graph density is $\frac{2m}{n(n-1)}$ and the maximum degree is d_{max} . The criteria of selecting these representative graphs are as follows. First, following [53], we do not select extremely easy or hard graphs, i.e., those graphs that can be solved within 5 seconds by all five solvers or cannot be solved within 3600 seconds by any solver when $k = 5$. Second, the representative graphs cover a wide range of sizes. Among the selected graphs, 10 small dense graphs (G1-G10) are synthetic graphs from **DIMACS-2 Graphs**, 10 medium graphs (G11-G20) with at most 10^6 vertices, and 10 large sparse graphs (G21-G30) with at least 10^6 vertices are real-world graphs from **Network Repository**. Third, most of the representative graphs have also been selected in previous studies [36, 53, 56].

6.1 Comparing with State-of-the-art Algorithms

Number of solved instances on two collections of graphs. We compare kPEX with four baselines by reporting the numbers of solved instances. The results for **Network Repository** are shown in Table 2 and Figure 4. We observe that kPEX outperforms all baselines for all tested values of k . For example, kPEX solves 12 instances more than the best baseline KPLEX for $k = 10$ within 3600 seconds. In addition, our kPEX is more stable than baselines when varying k . In contrast, there is an obvious drop in solved instances for kPlexT

¹<https://github.com/LijunChang/Maximum-kPlex>

²https://github.com/joey001/kplex_degen_gap

³<https://github.com/huajiang-ynu/ijcai23-kpx>

⁴<https://github.com/LijunChang/Maximum-kPlex-v2>

Table 2: Number of solved instances on Network Repository within 3600 seconds

k	kPEX (ours)	KPLEX	kPlexT	kPlexS	DiseMKP
2	567	564	559	559	542
3	564	553	557	553	527
5	565	557	554	547	516
10	564	552	537	549	495
15	559	548	507	547	452
20	559	546	471	539	439

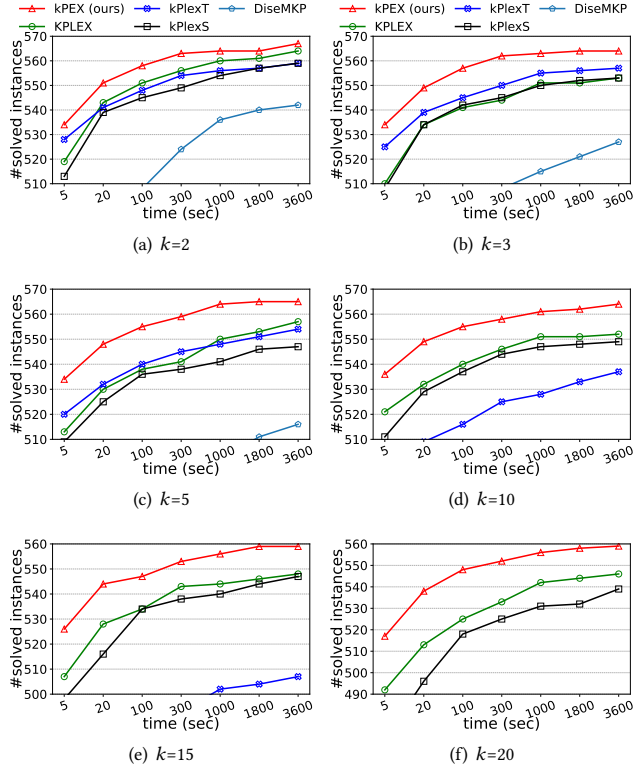


Figure 4: Number of solved instances on Network Repository (The lines corresponding to DiseMKP and kPlexT may not appear in the figures, as they are slow under certain settings and thus cannot reach the bottom lines within 3600 seconds.)

and DiseMKP as k increases from 2 to 20. This demonstrates the superiority of kPEX, which employs the AlTRB strategy (with novel reduction and bounding techniques) and efficient pre-processing methods. The results on the collection of **DIMACS-2 Graphs** are shown in Table 3 and Figure 5. kPEX outperforms all baselines by solving the most instances with 3600 seconds for all tested k values, e.g., kPEX solves 9 instances more than the second best solver KPLEX for $k = 5$. Besides, we note that kPEX is comparable with DiseMKP when $k = 2$. This is because the proposed reduction rules and upper bounding method are less effective for small values of k .

Table 3: Number of solved instances on DIMACS-2 within 3600 seconds

k	kPEX (ours)	KPLEX	kPlexT	kPlexS	DiseMKP
2	29	27	25	22	27
3	28	23	24	20	25
5	27	18	17	15	17
10	22	17	14	15	16
15	23	22	20	20	21
20	26	21	21	21	18

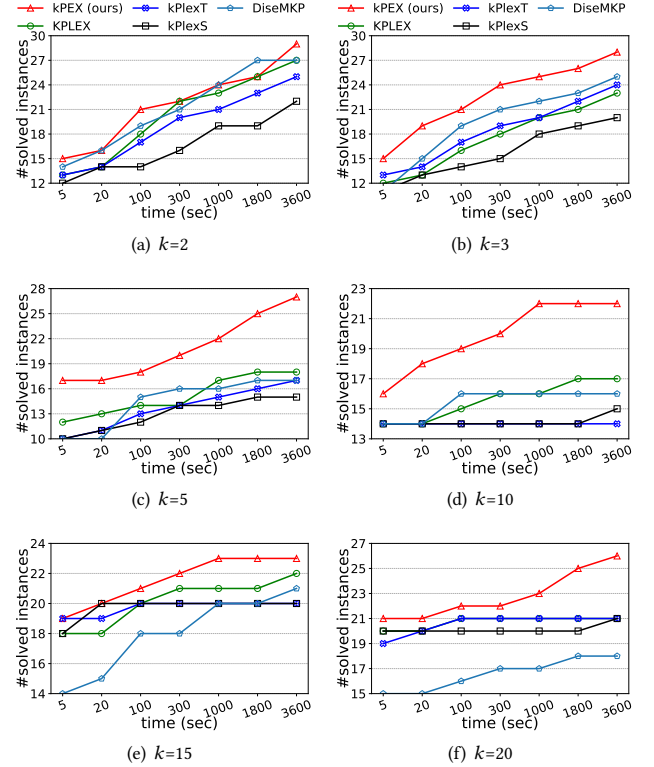


Figure 5: Number of solved instances on DIMACS-2

Running times on representative graphs. We report the running times of all algorithms on 30 representative graphs with $k = 5$ in Table 4. We observe that kPEX outperforms all baselines by achieving significant speedups on the majority graphs. For example, kPEX runs at least 5 times faster than KPLEX on 25 out of 30 graphs and at least 5 times faster than kPlexT on 21 out of 30 graphs. Moreover, there are 7 out of 30 graphs where kPEX runs at least 100 times faster than all baselines. Note that kPEX may exhibit slower performance compared to baselines on rare occasions. For instance, the baseline kPlexT runs faster than our kPEX on G23 with $k = 5$. The possible reasons include: 1) All algorithms rely on some heuristic procedures, e.g., the heuristic method for finding a large initial k -plex. The performance of these heuristic methods varies across different graphs and settings; 2) Compared with baselines, kPEX

Table 4: Running time in seconds of kPEX and state-of-the-arts on 30 graphs with $k = 5$

ID	kPEX (ours)	KPLEX	kPlexT	kPlexS	DiseMKP
G1	3.88	1154.76	120.63	187.88	55.56
G2	1360.94	OOT	OOT	OOT	OOT
G3	1527.41	OOT	OOT	OOT	OOT
G4	306.14	OOT	3160.85	OOT	OOT
G5	0.10	0.29	34.09	1356.45	0.16
G6	24.63	OOT	OOT	OOT	OOT
G7	433.86	OOT	OOT	OOT	OOT
G8	2.22	567.84	20.91	OOT	27.92
G9	2.82	411.26	OOT	OOT	OOT
G10	191.31	OOT	1339.31	OOT	1133.52
G11	3.97	1766.68	2318.35	OOT	OOT
G12	0.39	1.30	0.53	1.36	440.90
G13	55.29	OOT	OOT	OOT	OOT
G14	927.11	OOT	OOT	OOT	OOT
G15	21.17	OOT	OOT	OOT	OOT
G16	1.68	58.82	27.28	1655.40	900.80
G17	1.68	3022.10	123.86	OOT	OOT
G18	0.72	2804.46	1818.90	1099.39	OOT
G19	0.64	3.59	2.15	2.06	1695.05
G20	4.39	795.20	204.00	72.03	1347.27
G21	2.82	961.06	1515.40	OOT	OOT
G22	2.42	13.58	3.72	11.77	18.25
G23	13.16	136.61	4.80	11.39	OOT
G24	5.61	2979.37	3055.73	OOT	OOT
G25	3.80	92.05	203.26	OOT	OOT
G26	4.84	700.66	7.25	39.25	8.41
G27	2.60	14.52	3.50	15.39	51.94
G28	139.52	OOT	OOT	1703.49	OOT
G29	6.08	312.93	OOT	OOT	OOT
G30	593.26	OOT	OOT	OOT	OOT

Table 5: Running time in seconds of kPEX and its variants on 30 graphs with $k = 5$

ID	kPEX	kPEX-SeqRB	kPEX-CTCP	kPEX-EGo	kPEX-Degen
G1	3.88	18.66	3.89	3.91	3.92
G2	1360.94	OOT	1364.70	1371.95	1365.94
G3	1527.41	OOT	1525.00	1539.98	1538.84
G4	306.14	3220.90	305.95	307.76	307.69
G5	0.10	0.11	0.10	0.10	0.09
G6	24.63	557.07	24.63	33.37	57.68
G7	433.86	OOT	434.65	528.89	674.42
G8	2.22	18.54	2.22	2.19	2.20
G9	2.82	20.47	2.81	4.21	4.18
G10	191.31	3472.65	191.72	202.50	202.49
G11	3.97	23.51	4.63	4.19	3.87
G12	0.39	0.39	2.24	0.88	0.71
G13	55.29	555.36	59.38	97.39	103.23
G14	927.11	OOT	936.52	1113.52	1197.87
G15	21.17	128.27	35.76	29.85	24.84
G16	1.68	2.19	5.84	2.35	2.18
G17	1.68	8.49	8.27	3.88	1.90
G18	0.72	6.76	1.34	0.90	0.66
G19	0.64	0.66	17.41	8.36	93.70
G20	4.39	4.46	222.02	150.66	260.41
G21	2.82	8.03	3.55	3.68	3.32
G22	2.42	2.26	16.82	8.28	2.40
G23	13.16	12.62	1618.38	461.19	117.33
G24	5.61	15.80	6.97	9.83	9.03
G25	3.80	3.56	66.56	22.87	3.69
G26	4.84	4.77	4.64	5.61	4.54
G27	2.60	2.48	24.83	11.07	2.63
G28	139.52	139.96	OOT	OOT	OOT
G29	6.08	17.57	5.99	11.18	10.85
G30	593.26	OOT	1628.00	2094.29	2020.66

incorporates newly proposed reduction techniques, which may introduce additional time costs.

6.2 Effectiveness of Proposed Techniques

We compare the running time of kPEX with its variants:

- **kPEX-SeqRB**: kPEX replaces Al tRB with SeqRB(Section 4).
- **kPEX-CTCP**: kPEX replaces CF-CTCP with CTCP ([11]).
- **kPEX-EGo**: kPEX replaces KPHeur is with the existing heuristic method EGo-Degen in [12].
- **kPEX-Degen**: kPEX replaces KPHeur is with the existing heuristic method kPlex-Degen in [11].

In other words, kPEX-SeqRB is the version without Al tRB; kPEX-CTCP is the version without CF-CTCP; kPEX-EGo and kPEX-Degen are the versions without KPHeur is.

Effectiveness of Al tRB. We compare kPEX with kPEX-SeqRB and report the running times in Table 5. We observe that kPEX performs better than kPEX-SeqRB by achieving at least a 5 \times speedup on 12 out of 30 graphs and running up to 20 \times faster on G6. This indicates the effectiveness of Al tRB in narrowing down the search space. Besides, Al tRB contributes more speedups on synthetic graphs G1-G10 since the running time is dominated by the branch-reduction-and-bound stage on these graphs.

Effectiveness of CF-CTCP. We compare kPEX with kPEX-CTCP, and the running times are reported in Table 5. First, kPEX and kPEX-CTCP have similar performance on G1-G10 because the pre-processing techniques take little time (e.g., less than 1 second) on these synthetic graphs. Second, kPEX runs at least 5 times faster than kPEX-CTCP on 8 out of 20 real-world graphs. Moreover, CF-CTCP provides at least 50 \times speedup on G20 and G23. These results show the effectiveness of CF-CTCP on large sparse graphs.

Effectiveness of KPHeur is. We compare kPEX with its variants kPEX-EGo and kPEX-Degen (note that CF-CTCP is not replaced). The running times are shown in Table 5. We have the following observations. First, the running time of kPEX is less than that of both variants on the majority of graphs (i.e., on 23 out of 30 graphs). Then, kPEX runs at least 5 times faster than kPEX-EGo on 5 out of 30 graphs and faster than kPEX-Degen on 4 out of 30 graphs. In addition, kPEX runs at least 25 times faster than both kPEX-EGo and kPEX-Degen on G20 and G28. This shows that making more effort to finding a larger initial k -plex benefits kPEX by narrowing down the search space. Second, although kPEX may be slightly slower than the two variants, the extra time consumption is small and can be ignored compared to the total running time. For example, kPEX is 0.1 seconds slower than kPEX-Degen on G11 due to the extra computation, while the total running time of kPEX is 3.97

Table 6: Pre-processing time in seconds on 20 graphs with $k=5$ (lb denotes the size of the computed heuristic k -plex)

ID	kPEX		kPlexT		kPlexS		DiseMKP	
	time	lb	time	lb	time	lb	time	lb
G11	0.43	51	0.41	50	0.30	50	0.53	49
G12	0.39	73	0.58	69	1.30	34	1.71	34
G13	4.35	39	1.69	37	2.44	36	3.85	35
G14	9.24	70	5.01	67	3.08	64	6.05	66
G15	3.06	49	2.06	48	4.08	48	10.59	47
G16	1.29	27	0.82	26	1.84	14	3.85	14
G17	0.68	39	0.62	37	1.69	37	5.53	37
G18	0.14	87	0.34	87	0.27	87	0.51	87
G19	0.63	44	2.08	42	0.93	37	9.60	37
G20	3.48	21	33.60	19	32.56	10	118.24	10
G21	0.64	44	0.37	43	0.43	43	0.66	42
G22	2.42	34	4.51	32	15.46	26	15.55	27
G23	13.16	11	5.91	11	13.05	10	1708.75	10
G24	1.28	44	0.67	41	0.94	41	1.40	41
G25	3.14	77	7.90	76	34.82	76	45.58	77
G26	2.90	881	3.57	881	3.94	881	3.90	880
G27	2.60	37	4.79	35	18.02	32	21.06	33
G28	102.57	17	719.41	15	1210.59	12	OOT	-
G29	2.74	296	3.17	292	3.31	292	8.52	292
G30	102.73	65	134.74	62	252.29	17	1217.99	16

seconds, which means that the extra time consumption is negligible. Third, the performance of kPEX-EGo and kPEX-Degen is better than kPEX-SeqRB on G1-G10. This means that the variant of kPEX without AlTRB is slower than the variant without KPHeur_is. This indicates that AlTRB provides a greater performance boost than heuristic techniques on those graphs where branch-reduction-and-bound stage dominates the running time.

Effectiveness of KPHeur_is and CF-CTCP. We also compare the total pre-processing time and the size of the k -plex (i.e., lb) obtained by different heuristic methods in kPEX, kPlexT, kPlexS, and DiseMKP (note that KPLEX uses the same pre-processing method as kPlexS). The results are reported in Table 6. Note that we exclude the results on synthetic graphs G1-G10 since they have only hundreds of vertices and can be handled within 1 second by all methods. We have the following observations. First, kPEX consistently obtains the largest lb (or matches the largest obtained by others) while the pre-processing time remains comparable to other algorithms. Second, KPHeur_is outperforms the other pre-processing algorithms by obtaining a larger k -plex while costing much less time on G20 and G28. This also verifies the effectiveness of CF-CTCP and KPHeur_is.

7 RELATED WORK

Maximum k -plex search. The *maximum k -plex search* problem has garnered significant attention in social network analysis [42, 43] since the concept of k -plex was first proposed in [48]. Balasundaram et al. [3] showed the NP-hardness of the problem with any fixed k . Consequently, the major algorithmic design paradigm for exact solution is based on the *branch-reduction-and-bound* (BRB) framework [11, 12, 27, 35, 36, 53, 56, 63]. In particular, Xiao et al. [56] proposed a branching strategy, which improves theoretical time complexity from the trivial bound of $O^*(2^n)$ to $O^*(c^n)$ where $c < 2$ and O^* ignores polynomial factors. Later, Wang et al. [53] designed

KPLEX which is parameterized by the degeneracy gap (bounded empirically by $O(\log n)$). Very recently, Chang and Yao [12] proposed **kPlexT**, which improves the worst-case time complexity with newly proposed branching and reduction techniques. Additionally, several reduction and bounding techniques have been designed in the BRB framework to boost the practical performance. Gao et al. [27] developed reduction methods and a dynamic vertex selection strategy. Later, Zhou et al. [63] proposed a stronger reduction method and designed a coloring-based bounding method. Jiang et al. [36] designed a partition-based bounding method, and later in [35], their algorithm **DiseMKP** is equipped with a better upper bound. Chang et al. [11] designed an efficient algorithm **kPlexS** with a novel reduction method CTCP and a heuristic method. We note that the algorithms designed by Xiao et al. [56] and Chang and Yao [12] also work for the case when there is no requirement for the found k -plex to be of size at least $2k - 1$. We remark that existing works mainly focus on the BRB framework that conducts the reduction and the bounding sequentially, and our solution kPEX firstly adopts a new BRB framework that alternatively and iteratively conducts the reduction and the bounding.

Maximal k -plex enumeration. Another related problem is *maximal k -plex enumeration*, which aims to list all maximal k -plexes in the input graph; Here, a k -plex is *maximal* if it cannot be contained in other k -plexes. Many efficient algorithms are proposed for enumerating maximal k -plexes, including Bron-Kerbosch-based algorithms [17, 19, 22, 52, 54, 55] and reverse-search-based algorithms [6]. We remark that existing algorithms for enumerating maximal k -plexes can be utilized to solve the studied problem by listing all maximal k -plexes and then returning the largest one among them (note that the maximum k -plex is the maximal k -plex with largest number of vertices). However, the resulting solutions are not efficient due to the limited pruning and bounding techniques, as verified in [12].

Other cohesive subgraph models. k -plexes reduce to cliques when $k = 1$. There have been lines of work focusing on the maximum clique search and maximal clique enumeration problems [7, 8, 18, 24, 44, 45, 49, 50]. Further, the concept of k -plex is also explored in other kinds of graphs, e.g., bipartite graphs [13, 23, 40, 59, 60], directed graphs [30], temporal graphs [5], uncertain graphs [20], and so on. Besides k -plex, various cohesive subgraph models have been studied, including k -core [4, 15], k -truss [16, 33, 51], γ -quasi-clique [37, 46, 58, 61], k -defective clique [9, 14, 21, 28], densest subgraph [41, 57], and so on. For an overview on cohesive subgraph search, we refer to excellent books and surveys [10, 25, 26, 34, 39].

8 CONCLUSION

In this paper, we studied the maximum k -plex search problem. We proposed a new branch-reduction-and-bound method, called kPEX, which includes a new alternated reduction-and-bound process AlTRB. In addition, we also designed efficient pre-processing techniques for boosting the performance, which includes KPHeur_is for computing a large heuristic k -plex and CF-CTCP for efficiently removing unpromising vertices/edges. Extensive experiments on 664 graphs verified kPEX's superiority over state-of-the-art algorithms. In the future, we will explore the possibility of adapting kPEX to mining other cohesive subgraphs.

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A ADDITIONAL DESCRIPTIONS OF CF-CTCP

A.1 Time Complexity of CF-CTCP

Before analyzing the time complexity of CF-CTCP (Algorithm 4), we first prove the following lemma.

LEMMA A.1. *Given a graph $G = (V, E)$, we have*

$$\sum_{(u,v) \in E} \min(d_G(u), d_G(v)) \leq 2m \times \delta(G).$$

PROOF. Assume that vertices v_1, v_2, \dots, v_n in G are sorted according to the degeneracy order, indicating that $|N_G^+(v_i)| = |N_G(v_i) \cap \{v_{i+1}, v_{i+2}, \dots, v_n\}| \leq \delta(G)$. Thus we have

$$\begin{aligned} & \sum_{(u,v) \in E} \min(d_G(u), d_G(v)) \\ &= \sum_{v_i \in V} \sum_{v_j \in N_G^+(v_i)} \min(d_G(v_i), d_G(v_j)) \\ &\leq \sum_{v_i \in V} \sum_{v_j \in N_G^+(v_i)} d_G(v_i) \leq \sum_{v_i \in V} d_G(v_i) \times \delta(G) = 2m \times \delta(G). \end{aligned}$$

□

We can derive from Lemma A.1 that

$$O\left(\sum_{(u,v) \in E} \min(d_G(u), d_G(v))\right) = O(m \times \delta(G)).$$

Now we are ready to prove the total time complexity of CF-CTCP (Lemma 5.3).

PROOF. Note that we invoke CF-CTCP only when $Q_v \neq \emptyset$ or $lb_changed = true$ as in CTCP [11]. First, for the first invocation, Line 6 of Algorithm 4 computes the common neighbors $\Delta(u, v)$ for each edge (u, v) , and the time complexity is

$$O\left(\sum_{(u,v) \in E} \min(d_G(u), d_G(v))\right) = O(m \times \delta(G)),$$

according to Lemma A.1. Second, the total time consumption of core pruning is $O(m)$ [4] and the total time cost of Procedure RemoveEdge is also $O(m)$ since we can implement Line 21 for at most m times. Third, for all invocations, there are at most $\delta(G)$ times when $lb_changed = true$ since $k \leq lb = |S^*| \leq \delta(G) + k$ (S^* denoting the largest k -plex seen so far), which indicates that we will perform Lines 4-8 at most $\delta(G)$ times. Thus the total time complexity of Lines 1-8 is $O(m \times \delta(G))$. We next consider Lines 9-20. We will pop at most m edges, and for each edge, we need to find all the triangles that it participates in, which can be done in $O(\sum_{(u,v) \in E} \min(d_G(u), d_G(v))) = O(m \times \delta(G))$. Therefore, the total time complexity of all invocations to CF-CTCP is $O(m \times \delta(G))$, which completes our proof. □

A.2 An Implementation of CF-CTCP with $O(m)$ Memory

A direct implementation of CF-CTCP requires storing the common neighbors $\Delta(\cdot, \cdot)$ for all edges, which needs $O(m \times \delta(G))$ memory. In the following, we propose a novel implementation that requires only $O(m)$ memory without changing the time complexity of CF-CTCP. In particular, we need three auxiliary arrays A_1, A_2 , and A_3 , each of length m , to store additional information for each edge: 1) array A_1

records the number of triangles, 2) array A_2 records the timestamp (e.g., system time) when the triangle count is computed in Line 6, and 3) array A_3 records the timestamp (e.g., system time) when an edge is removed in Lines 1, 21 and 22. Based on these three arrays, we correspondingly modify Algorithm 4 as follows. First, we only record $|\Delta(u, v)|$ using A_1 instead of storing the whole vertex set $\Delta(u, v)$ in Line 6. The correspond triangle count in A_1 is decreased by 1 when CF-CTCP modifies $\Delta(\cdot, \cdot)$ in Lines 13 and 18. Second, when we traverse all triangles that edge (u, v) belongs to in Line 12, we enumerate such a vertex w that satisfies: 1) both (u, w) and (v, w) are in $E \cup Q_e$, i.e., (u, v, w) forms a triangle; 2) the timestamp of computing the triangle count for edge (u, w) is before the timestamp of removing edge (u, v) using arrays A_2 and A_3 , i.e., when we compute $|\Delta(u, w)|$ in Line 6, edge (u, v) has not yet been removed. The modification to Line 17 follows the same fashion as Line 12. Finally, it is easy to verify the correctness of the above modification of CF-CTCP with $O(m)$ memory usage.

B ADDITIONAL EXPERIMENTAL RESULTS

We provide additional experimental results for $k = 2, 3, 10, 15, 20$.

B.1 Comparing with State-of-the-art Algorithms

Running times on representative graphs. We report the running times of all algorithms on 30 representative graphs with $k = 2, 3, 10, 15$, and 20 in Tables 7, 8, 9, 10, and 11, respectively. We observe that kPEX outperforms all baselines by achieving significant speedups on the majority graphs. For example, kPEX runs at least 5 times faster than KPLEX on 23 out of 30 graphs and at least 5 times faster than kPlExT on 20 out of 30 graphs when $k = 3$.

B.2 Effectiveness of Proposed Techniques

Effectiveness of AltrB. We compare kPEX with kPEX-SeqRB and report the running times for $k=2, 3, 10, 15$, and 20 in Tables 12, 13, 14, 15, and 16, respectively. We observe that kPEX performs better than kPEX-SeqRB at most times, and AltrB can bring at least 60× speedup on G5 when $k=10$. In addition, we observe that the gap between kPEX and kPEX-AltrB narrows when $k \geq 15$. A possible reason may be that finding a larger heuristic k -plex (i.e., KPHeur is) is more important than AltrB for large values of k .

Effectiveness of CF-CTCP. We compare kPEX with kPEX-CTCP, and the running times for $k=2, 3, 10, 15$, and 20 are reported in Tables 12, 13, 14, 15, and 16, respectively. First, kPEX and kPEX-CTCP still have similar performance on G1-G10 because the pre-processing techniques take little time on these small synthetic graphs. Second, kPEX runs stably at least 50 times faster than kPEX-CTCP on G23 for all tested values of k . These results show the effectiveness of CF-CTCP on large sparse graphs.

Effectiveness of KPHeur is. We compare kPEX with its variants kPEX-EGo and kPEX-Degen (note that CF-CTCP is not replaced). The running times for $k=2, 3, 10, 15$, and 20 are shown in Tables 12, 13, 14, 15, and 16, respectively. We have the following observations. First, the running time of kPEX is less than that of both variants on the majority of graphs. Then, kPEX runs up to three orders of magnitude faster than both kPEX-EGo and kPEX-Degen

Table 7: Running time in seconds of kPEX and state-of-the-arts on 30 graphs with $k = 2$

ID	kPEX (ours)	KPLEX	kPlexT	kPlexS	DiseMKP
G1	0.95	1.77	1.56	2.73	1.23
G2	1982.78	1847.75	OOT	OOT	OOT
G3	51.67	105.06	128.29	178.26	21.63
G4	1.35	9.46	8.02	18.98	2.08
G5	77.13	27.08	OOT	OOT	OOT
G6	OOT	OOT	OOT	OOT	OOT
G7	OOT	OOT	OOT	OOT	OOT
G8	0.33	2.60	2.47	18.91	0.39
G9	39.77	86.99	87.66	637.09	497.46
G10	4.07	33.39	30.03	479.11	3.30
G11	20.89	127.32	121.80	1009.49	OOT
G12	0.54	1.45	0.70	1.56	13.80
G13	289.18	2505.66	2797.56	OOT	2401.74
G14	3173.74	OOT	OOT	OOT	OOT
G15	142.05	1420.48	1627.85	OOT	OOT
G16	1.96	7.69	8.05	31.33	7.98
G17	6.30	39.70	51.66	220.65	101.81
G18	2.28	8.16	71.92	61.06	OOT
G19	8.71	11.07	28.43	9.39	OOT
G20	5.68	23.69	77.29	19.64	328.39
G21	13.00	118.76	123.49	942.62	337.35
G22	3.36	16.27	3.79	15.74	21.88
G23	8.66	10.12	3.87	10.54	1035.93
G24	24.66	258.56	245.02	2345.22	550.65
G25	13.90	79.16	77.10	277.82	OOT
G26	11.59	41.78	5.98	107.83	232.15
G27	3.18	19.32	4.39	17.14	26.33
G28	145.64	849.77	OOT	1347.36	OOT
G29	6.02	7.71	2563.26	535.31	OOT
G30	147.23	514.32	721.71	1492.62	OOT

on G19 when $k=20$. This shows that making more effort to finding a larger initial k -plex benefits kPEX by narrowing down the search space.

Effectiveness of KPHeuris and CF-CTCP. We also compare the total pre-processing time and the size of the k -plex (i.e., lb) obtained by different heuristic methods in kPEX, kPlexT, kPlexS, and DiseMKP (note that KPLEX uses the same pre-processing method as kPlexS). The results for $k=2, 3, 10, 15$, and 20 are reported in Tables 17, 18, 19, 20, and 21, respectively. Note that we exclude the results on synthetic graphs G1-G10 since they have only hundreds of vertices and can be handled within 1 second by all methods. We have the following observations. First, kPEX obtains the largest lb (or matches the largest obtained by others) at most time while the pre-processing time remains comparable to other algorithms. Second, KPHeuris outperforms the other pre-processing algorithms by obtaining a larger k -plex while costing much less time on G20 and G22 for all tested values of k . This also verifies the effectiveness of CF-CTCP and KPHeuris.

Table 8: Running time in seconds of kPEX and state-of-the-arts on 30 graphs with $k = 3$

ID	kPEX (ours)	KPLEX	kPlexT	kPlexS	DiseMKP
G1	5.38	32.14	23.32	22.39	7.30
G2	OOT	OOT	OOT	OOT	OOT
G3	60.32	1112.70	2071.59	1461.59	22.36
G4	9.17	269.58	69.18	705.19	10.98
G5	0.10	0.46	1.22	15.65	0.08
G6	2.62	28.08	1552.41	OOT	49.09
G7	OOT	OOT	OOT	OOT	OOT
G8	0.16	73.99	2.78	322.77	1.02
G9	21.19	528.72	953.50	3592.43	2197.53
G10	15.04	2622.38	164.85	OOT	10.05
G11	6.55	OOT	374.45	OOT	OOT
G12	0.41	1.46	0.61	1.54	15.69
G13	163.63	OOT	OOT	OOT	OOT
G14	OOT	OOT	OOT	OOT	OOT
G15	46.87	OOT	2702.41	OOT	OOT
G16	1.46	167.99	7.66	184.59	20.50
G17	2.14	987.10	35.01	926.33	245.60
G18	1.73	577.38	292.77	314.95	OOT
G19	0.98	3.36	0.85	1.34	2921.83
G20	4.28	102.50	26.06	25.77	317.76
G21	8.88	OOT	349.83	OOT	1712.85
G22	3.23	15.34	3.63	14.41	21.18
G23	13.17	150.23	4.55	10.65	1761.28
G24	15.21	OOT	643.10	OOT	OOT
G25	6.92	2877.27	181.87	1930.20	OOT
G26	7.07	368.98	10.68	251.34	50.96
G27	3.01	17.59	3.73	17.10	27.71
G28	117.66	OOT	OOT	1163.18	OOT
G29	115.24	825.59	OOT	OOT	OOT
G30	200.21	OOT	OOT	OOT	OOT

Table 9: Running time in seconds of kPEX and state-of-the-arts on 30 graphs with $k = 10$

ID	kPEX (ours)	KPLEX	kPlexT	kPlexS	DiseMKP
G1	462.04	OOT	OOT	3142.77	OOT
G2	38.55	OOT	OOT	OOT	OOT
G3	OOT	OOT	OOT	OOT	OOT
G4	OOT	OOT	OOT	OOT	OOT
G5	14.63	OOT	OOT	OOT	OOT
G6	OOT	OOT	OOT	OOT	OOT
G7	OOT	OOT	OOT	OOT	OOT
G8	OOT	OOT	OOT	OOT	OOT
G9	309.46	OOT	OOT	OOT	OOT
G10	OOT	OOT	OOT	OOT	OOT
G11	6.00	OOT	OOT	OOT	OOT
G12	0.50	1.12	0.58	1.40	OOT
G13	859.77	OOT	OOT	OOT	OOT
G14	3201.23	OOT	OOT	OOT	OOT
G15	23.07	OOT	OOT	OOT	OOT
G16	2.01	23.08	OOT	293.97	OOT
G17	2.03	314.96	OOT	OOT	OOT
G18	0.51	2679.72	OOT	1017.07	OOT
G19	0.29	3.04	5.21	2.65	1388.96
G20	5.17	409.33	OOT	31.76	OOT
G21	20.86	OOT	OOT	OOT	OOT
G22	2.07	11.16	2.87	10.38	18.09
G23	13.47	163.83	4.81	12.85	OOT
G24	34.49	OOT	OOT	OOT	OOT
G25	2.81	34.57	OOT	587.10	OOT
G26	1.10	3.61	3.73	3.90	4.04
G27	2.47	15.41	3.37	15.06	OOT
G28	263.01	OOT	OOT	OOT	OOT
G29	2.89	101.16	OOT	OOT	11.28
G30	OOT	OOT	OOT	OOT	OOT

Table 10: Running time in seconds of kPEX and state-of-the-arts on 30 graphs with $k = 15$

ID	kPEX (ours)	KPLEX	kPlexT	kPlexS	DiseMKP
G1	7.84	44.29	35.84	17.20	418.76
G2	0.06	34.53	3.58	8.27	22.95
G3	OOT	OOT	OOT	OOT	OOT
G4	OOT	OOT	OOT	OOT	OOT
G5	OOT	OOT	OOT	OOT	OOT
G6	OOT	OOT	OOT	OOT	OOT
G7	OOT	OOT	OOT	OOT	OOT
G8	OOT	OOT	OOT	OOT	OOT
G9	115.99	OOT	OOT	OOT	OOT
G10	OOT	OOT	OOT	OOT	OOT
G11	18.38	OOT	OOT	OOT	OOT
G12	0.47	1.08	0.57	1.07	OOT
G13	1338.78	OOT	OOT	OOT	OOT
G14	OOT	OOT	OOT	OOT	OOT
G15	141.05	OOT	OOT	OOT	OOT
G16	1.78	2.02	OOT	2.63	OOT
G17	6.66	14.37	OOT	321.01	OOT
G18	0.12	0.33	1144.04	0.28	2621.37
G19	0.39	4.43	3.07	4.23	OOT
G20	594.86	194.44	OOT	139.13	OOT
G21	122.70	OOT	OOT	OOT	OOT
G22	1.68	8.63	33.25	9.01	OOT
G23	13.21	132.50	4.81	11.67	OOT
G24	288.71	OOT	OOT	OOT	OOT
G25	5.74	168.11	OOT	OOT	OOT
G26	1.12	3.63	3.57	3.88	4.02
G27	1.79	11.92	OOT	9.44	OOT
G28	OOT	OOT	OOT	OOT	OOT
G29	2.46	4.08	23.15	3.69	36.37
G30	OOT	OOT	OOT	OOT	OOT

Table 11: Running time in seconds of kPEX and state-of-the-arts on 30 graphs with $k = 20$

ID	kPEX (ours)	KPLEX	kPlexT	kPlexS	DiseMKP
G1	0.00	0.00	0.00	0.00	0.00
G2	0.00	0.00	0.00	0.00	0.00
G3	OOT	OOT	OOT	OOT	OOT
G4	OOT	OOT	OOT	OOT	OOT
G5	OOT	OOT	OOT	OOT	OOT
G6	OOT	OOT	OOT	OOT	OOT
G7	OOT	OOT	OOT	OOT	OOT
G8	OOT	OOT	OOT	OOT	OOT
G9	36.06	OOT	OOT	OOT	OOT
G10	OOT	OOT	OOT	OOT	OOT
G11	35.88	OOT	OOT	OOT	OOT
G12	0.30	1.08	0.53	0.99	OOT
G13	1022.03	OOT	OOT	OOT	OOT
G14	3067.04	OOT	OOT	OOT	OOT
G15	409.76	OOT	OOT	OOT	OOT
G16	9.53	311.28	OOT	2457.82	OOT
G17	47.76	41.72	OOT	1801.99	OOT
G18	0.19	0.83	OOT	0.99	OOT
G19	0.44	4.50	2.61	5.01	OOT
G20	114.36	1398.04	OOT	59.27	OOT
G21	409.39	OOT	OOT	OOT	OOT
G22	1.77	7.64	OOT	8.22	OOT
G23	12.89	130.41	5.05	28.77	OOT
G24	1045.59	OOT	OOT	OOT	OOT
G25	10.56	334.13	OOT	OOT	OOT
G26	1.05	3.80	3.42	3.67	2.70
G27	2.10	11.52	OOT	11.18	OOT
G28	OOT	OOT	OOT	OOT	OOT
G29	2.47	2.99	2.95	3.28	12.19
G30	OOT	OOT	OOT	OOT	OOT

Table 12: Running time in seconds of kPEX and its variants on 30 graphs with $k = 2$

ID	kPEX	kPEX-SeqRB	kPEX-CTCP	kPEX-EGo	kPEX-Degen
G1	0.95	1.23	0.95	0.95	0.95
G2	1982.78	2021.27	1988.00	1992.98	1994.67
G3	51.67	75.39	51.66	52.01	52.18
G4	1.35	2.90	1.34	1.36	1.36
G5	77.13	118.45	77.47	77.41	77.25
G6	OOT	OOT	OOT	OOT	OOT
G7	OOT	OOT	OOT	OOT	OOT
G8	0.33	0.29	0.33	0.29	0.31
G9	39.77	50.32	39.81	39.97	41.21
G10	4.07	5.16	4.05	3.81	3.88
G11	20.89	33.06	21.47	23.02	25.65
G12	0.54	0.55	2.41	0.96	1.12
G13	289.18	407.48	291.83	345.93	343.31
G14	3173.74	OOT	3186.76	OOT	OOT
G15	142.05	189.20	158.75	172.91	277.23
G16	1.96	2.00	6.23	3.34	3.33
G17	6.30	7.06	13.16	9.55	7.25
G18	2.28	2.50	2.89	2.39	2.56
G19	8.71	8.85	33.79	16.26	OOT
G20	5.68	4.91	245.87	190.31	316.79
G21	13.00	22.35	13.87	20.14	19.76
G22	3.36	3.20	22.43	11.40	3.43
G23	8.66	8.11	1418.09	469.20	8.54
G24	24.66	41.83	26.23	24.27	36.39
G25	13.90	17.56	86.75	35.87	13.80
G26	11.59	21.19	11.03	12.31	11.56
G27	3.18	3.31	30.24	13.21	3.58
G28	145.64	137.51	OOT	OOT	OOT
G29	6.02	7.65	6.46	10.13	10.28
G30	147.23	179.48	1453.76	1437.70	251.08

Table 13: Running time in seconds of kPEX and its variants on 30 graphs with $k = 3$

ID	kPEX	kPEX-SeqRB	kPEX-CTCP	kPEX-EGo	kPEX-Degen
G1	5.38	16.99	5.40	5.41	5.42
G2	OOT	OOT	OOT	OOT	OOT
G3	60.32	530.53	60.48	60.65	60.88
G4	9.17	59.21	9.20	9.53	9.52
G5	0.10	0.11	0.10	0.10	0.09
G6	2.62	21.44	2.62	2.85	3.18
G7	OOT	OOT	OOT	OOT	OOT
G8	0.16	0.18	0.15	0.17	0.17
G9	21.19	149.41	21.22	25.29	28.53
G10	15.04	82.46	15.03	14.80	15.20
G11	6.55	33.43	7.15	11.34	12.64
G12	0.41	0.41	2.29	0.92	0.91
G13	163.63	985.84	167.20	195.70	208.35
G14	OOT	OOT	OOT	OOT	OOT
G15	46.87	204.47	62.51	63.44	57.47
G16	1.46	1.68	5.91	2.26	1.87
G17	2.14	3.31	8.94	4.53	2.76
G18	1.73	3.96	2.30	2.36	2.49
G19	0.98	1.00	24.22	8.57	49.43
G20	4.28	4.40	226.14	171.87	271.25
G21	8.88	48.31	9.66	9.64	9.53
G22	3.23	2.81	20.21	9.99	3.34
G23	13.17	12.77	1617.71	468.21	118.89
G24	15.21	80.36	16.67	16.34	16.61
G25	6.92	13.36	76.74	28.93	7.86
G26	7.07	7.14	6.34	7.57	6.75
G27	3.01	2.88	27.60	12.35	3.18
G28	117.66	116.05	OOT	OOT	3271.82
G29	115.24	455.48	115.38	113.24	118.54
G30	200.21	833.98	1502.80	1356.16	353.15

Table 14: Running time in seconds of kPEX and its variants on 30 graphs with $k = 10$

ID	kPEX	kPEX-SeqRB	kPEX-CTCP	kPEX-EGo	kPEX-Degen
G1	462.04	1238.20	463.16	464.53	464.13
G2	38.55	50.23	38.79	38.71	38.58
G3	OOT	OOT	OOT	OOT	OOT
G4	OOT	OOT	OOT	OOT	OOT
G5	14.63	912.36	14.65	14.72	14.67
G6	OOT	OOT	OOT	OOT	OOT
G7	OOT	OOT	OOT	OOT	OOT
G8	OOT	OOT	OOT	OOT	OOT
G9	309.46	837.83	310.32	625.51	623.68
G10	OOT	OOT	OOT	OOT	OOT
G11	6.00	26.39	6.66	17.61	17.30
G12	0.50	0.52	2.22	0.83	0.60
G13	859.77	OOT	866.51	1206.11	1200.79
G14	3201.23	OOT	3207.34	3366.27	3346.56
G15	23.07	234.55	35.02	35.29	31.29
G16	2.01	3.71	6.10	4.78	4.75
G17	2.03	13.22	7.91	6.87	5.07
G18	0.51	21.17	1.09	3.77	3.57
G19	0.29	0.29	2.81	44.58	OOT
G20	5.17	7.01	204.72	108.26	225.67
G21	20.86	125.24	21.52	50.77	50.31
G22	2.07	1.97	15.59	7.08	2.05
G23	13.47	12.81	1613.00	462.58	116.71
G24	34.49	195.37	35.84	52.85	51.96
G25	2.81	2.73	58.16	17.78	2.91
G26	1.10	1.09	2.87	2.55	1.48
G27	2.47	2.33	23.40	9.59	2.38
G28	263.01	261.12	OOT	OOT	OOT
G29	2.89	3.18	2.86	2.76	2.58
G30	OOT	OOT	OOT	OOT	OOT

Table 15: Running time in seconds of kPEX and its variants on 30 graphs with $k = 15$

ID	kPEX	kPEX-SeqRB	kPEX-CTCP	kPEX-EGo	kPEX-Degen
G1	7.84	10.63	7.85	7.72	7.94
G2	0.06	0.06	0.05	0.06	0.05
G3	OOT	OOT	OOT	OOT	OOT
G4	OOT	OOT	OOT	OOT	OOT
G5	OOT	OOT	OOT	OOT	OOT
G6	OOT	OOT	OOT	OOT	OOT
G7	OOT	OOT	OOT	OOT	OOT
G8	OOT	OOT	OOT	OOT	OOT
G9	115.99	120.17	115.87	113.30	118.43
G10	OOT	OOT	OOT	OOT	OOT
G11	18.38	19.24	18.37	71.93	74.15
G12	0.47	0.47	2.11	0.78	0.62
G13	1338.78	1497.62	1329.66	2796.92	2867.16
G14	OOT	OOT	OOT	OOT	OOT
G15	141.05	167.43	148.50	170.58	173.08
G16	1.78	1.86	5.55	2.48	2.98
G17	6.66	7.52	12.19	7.16	5.70
G18	0.12	0.13	0.67	0.34	0.18
G19	0.39	0.40	0.70	728.27	OOT
G20	594.86	1240.20	785.65	943.65	1486.61
G21	122.70	131.19	121.98	225.87	233.15
G22	1.68	1.68	12.05	5.75	1.66
G23	13.21	13.15	1605.10	458.65	116.44
G24	288.71	316.40	286.53	281.00	293.41
G25	5.74	5.74	55.64	19.29	5.72
G26	1.12	1.10	2.72	2.49	1.48
G27	1.79	1.75	17.87	7.54	1.81
G28	OOT	OOT	OOT	OOT	OOT
G29	2.46	2.57	2.54	2.15	2.10
G30	OOT	OOT	OOT	OOT	OOT

Table 16: Running time in seconds of kPEX and its variants on 30 graphs with $k = 20$

ID	kPEX	kPEX-SeqRB	kPEX-CTCP	kPEX-EGo	kPEX-Degen
G1	0.00	0.00	0.00	0.00	0.00
G2	0.00	0.00	0.00	0.00	0.00
G3	OOT	OOT	OOT	OOT	OOT
G4	OOT	OOT	OOT	OOT	OOT
G5	OOT	OOT	OOT	OOT	OOT
G6	OOT	OOT	OOT	OOT	OOT
G7	OOT	OOT	OOT	OOT	OOT
G8	OOT	OOT	OOT	OOT	OOT
G9	36.06	35.42	35.87	71.18	74.99
G10	OOT	OOT	OOT	OOT	OOT
G11	35.88	35.96	35.82	58.23	59.62
G12	0.30	0.30	1.88	0.81	0.56
G13	1022.03	1010.12	1020.25	OOT	OOT
G14	3067.04	3450.98	3055.50	OOT	OOT
G15	409.76	416.74	412.44	1627.16	1688.80
G16	9.53	10.78	10.71	9.37	24.72
G17	47.76	58.52	52.69	48.09	49.33
G18	0.19	0.19	0.67	0.35	0.21
G19	0.44	0.45	0.57	OOT	OOT
G20	114.36	122.56	274.94	625.36	OOT
G21	409.39	408.68	407.72	572.13	600.41
G22	1.77	1.76	12.48	5.96	2.10
G23	12.89	12.65	1602.86	452.09	116.28
G24	1045.59	1060.97	1040.63	1134.76	1191.44
G25	10.56	10.50	58.15	24.76	11.29
G26	1.05	1.10	2.36	2.34	1.39
G27	2.10	2.10	16.94	7.54	2.17
G28	OOT	OOT	OOT	OOT	OOT
G29	2.47	2.47	2.45	1.91	1.84
G30	OOT	OOT	OOT	OOT	878.37

Table 17: Pre-processing time in seconds on 20 graphs with $k=2$ (lb denotes the size of the computed heuristic k -plex)

ID	kPEX		kPlexT		kPlexS		DiseMKP	
	time	lb	time	lb	time	lb	time	lb
G11	1.08	37	0.46	37	0.27	34	0.54	35
G12	0.52	63	0.77	62	1.32	21	1.78	23
G13	9.08	26	1.64	25	2.21	25	3.70	23
G14	25.60	54	5.17	52	2.91	51	6.12	52
G15	7.72	36	2.80	36	4.09	32	13.11	34
G16	1.16	19	0.91	17	1.96	9	4.07	9
G17	1.17	28	0.72	27	1.86	27	6.02	26
G18	0.33	70	0.42	70	0.29	69	0.55	70
G19	4.48	35	1.05	34	0.79	30	15.29	31
G20	4.02	15	4.10	14	9.06	10	141.38	4
G21	1.01	32	0.41	30	0.44	30	0.69	28
G22	3.36	31	5.58	31	20.15	17	21.32	18
G23	8.66	6	4.92	5	13.76	5	990.37	5
G24	1.75	32	0.82	32	0.98	30	1.58	30
G25	4.37	60	8.81	60	37.18	60	56.61	59
G26	3.01	872	3.57	872	3.76	872	3.20	872
G27	3.12	27	5.27	27	20.55	24	24.79	24
G28	102.61	11	72.20	10	201.25	9	OOT	-
G29	2.68	273	3.28	274	3.14	271	9.38	272
G30	108.37	52	144.45	51	287.50	10	1688.08	10

Table 18: Pre-processing time in seconds on 20 graphs with $k=3$ (lb denotes the size of the computed heuristic k -plex)

ID	kPEX		kPlexT		kPlexS		DiseMKP	
	time	lb	time	lb	time	lb	time	lb
G11	0.82	44	0.45	42	0.28	40	0.55	41
G12	0.40	67	0.70	66	1.38	26	1.72	27
G13	7.58	30	1.89	29	2.28	28	3.72	27
G14	11.84	58	5.23	58	3.23	56	6.17	56
G15	4.70	41	2.24	40	4.22	40	11.86	39
G16	1.10	22	0.78	21	1.76	11	4.15	10
G17	0.84	33	0.70	32	1.79	31	5.91	29
G18	0.45	77	0.40	75	0.28	74	0.54	74
G19	0.97	39	0.98	39	0.74	34	14.34	34
G20	3.39	17	4.49	16	8.21	10	133.31	6
G21	0.99	35	0.44	34	0.45	33	0.68	33
G22	3.23	32	5.63	31	20.64	20	22.30	18
G23	13.17	7	5.79	7	11.91	6	1723.31	6
G24	1.70	35	0.84	34	1.00	31	1.58	31
G25	3.83	66	9.69	66	35.19	64	52.06	65
G26	3.24	875	3.51	875	3.57	875	3.32	875
G27	3.01	32	5.16	32	19.15	27	24.36	28
G28	77.67	13	152.26	12	320.38	9	OOT	-
G29	2.53	280	3.12	280	3.47	280	9.15	279
G30	118.35	57	143.21	55	259.23	12	1481.51	12

Table 19: Pre-processing time in seconds on 20 graphs with $k=10$ (lb denotes the size of the computed heuristic k -plex)

ID	kPEX		kPlexT		kPlexS		DiseMKP	
	time	lb	time	lb	time	lb	time	lb
G11	0.25	67	0.37	65	0.28	65	0.49	67
G12	0.50	82	0.64	74	1.18	45	1.58	45
G13	2.55	54	1.42	52	2.55	52	3.91	54
G14	4.58	90	4.55	88	3.39	88	6.06	87
G15	1.84	65	1.85	64	3.73	64	8.75	62
G16	0.73	40	0.77	36	1.71	25	3.52	24
G17	0.64	53	0.59	48	1.66	48	5.04	49
G18	0.09	102	0.29	101	0.23	101	0.46	101
G19	0.28	53	5.25	47	1.18	44	1.75	44
G20	3.99	31	33.64	29	31.91	20	79.90	20
G21	0.53	57	0.55	54	0.42	54	0.61	54
G22	2.07	44	3.94	38	13.23	34	13.02	35
G23	13.47	21	5.90	21	13.03	20	1688.87	20
G24	0.89	57	0.63	55	0.95	55	1.26	54
G25	2.43	92	6.75	91	30.65	91	40.20	91
G26	1.10	891	4.33	891	4.75	891	4.20	891
G27	2.47	46	4.62	45	17.20	40	19.06	41
G28	241.44	27	1808.29	22	3247.72	21	OOT	-
G29	2.44	316	3.25	316	3.33	316	9.32	315
G30	99.21	82	124.70	77	229.23	26	887.60	26

Table 20: Pre-processing time in seconds on 20 graphs with $k=15$ (lb denotes the size of the computed heuristic k -plex)

ID	kPEX		kPlexT		kPlexS		DiseMKP	
	time	lb	time	lb	time	lb	time	lb
G11	0.20	79	0.34	77	0.26	77	0.45	77
G12	0.46	89	0.62	78	1.16	54	1.47	55
G13	2.02	67	1.37	63	2.42	63	3.86	65
G14	2.85	108	4.38	107	3.41	107	5.90	107
G15	1.46	77	1.39	76	3.45	76	7.58	74
G16	0.86	51	0.83	47	1.55	33	3.01	33
G17	0.45	64	0.56	59	1.57	59	4.76	57
G18	0.06	116	0.25	115	0.21	115	0.41	115
G19	0.38	59	2.35	50	1.22	49	0.42	49
G20	4.80	41	28.86	37	28.31	30	59.46	30
G21	0.44	69	0.52	67	0.40	67	0.59	68
G22	1.68	49	3.49	42	10.83	42	10.67	42
G23	13.21	31	5.81	31	12.79	30	1713.30	30
G24	0.66	69	0.59	69	0.82	69	1.15	66
G25	2.26	101	6.66	101	29.13	101	37.67	101
G26	1.12	900	4.22	900	4.24	900	3.91	900
G27	1.78	53	3.71	51	13.72	51	17.02	50
G28	469.30	36	1307.77	31	3177.53	31	OOT	-
G29	2.32	332	3.62	332	3.72	332	9.04	331
G30	86.23	98	111.59	92	215.72	36	659.92	36

Table 21: Pre-processing time in seconds on 20 graphs with $k=20$ (lb denotes the size of the computed heuristic k -plex)

ID	kPEX		kPlexT		kPlexS		DiseMKP	
	time	lb	time	lb	time	lb	time	lb
G11	0.17	89	0.31	88	0.24	88	0.41	87
G12	0.30	96	0.58	78	1.08	64	1.42	63
G13	2.36	79	1.33	76	2.40	76	3.60	77
G14	2.18	123	4.21	122	3.25	122	5.70	121
G15	1.33	88	1.67	85	3.35	85	6.55	86
G16	2.75	59	1.31	51	0.80	40	1.41	40
G17	0.46	74	0.57	67	1.51	67	4.54	67
G18	0.06	124	0.24	123	0.20	123	0.39	123
G19	0.33	64	1.08	54	1.01	54	0.31	54
G20	5.30	50	31.26	44	25.04	40	45.51	40
G21	0.64	79	0.49	76	0.39	76	0.57	78
G22	1.76	55	5.07	47	11.81	47	11.71	48
G23	12.89	41	5.90	41	13.27	40	1695.05	40
G24	0.77	78	0.58	77	0.76	77	1.08	75
G25	2.22	111	6.75	110	25.55	110	35.31	110
G26	1.05	910	4.25	910	4.29	910	2.94	910
G27	1.76	60	3.96	58	12.15	58	14.96	58
G28	1986.00	45	779.28	40	2487.34	40	947.97	40
G29	2.38	343	3.33	343	3.62	343	8.70	342
G30	76.54	112	98.14	104	200.47	46	554.62	45